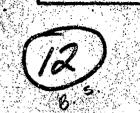
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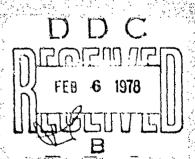
A VERSATILE USER-ORIENTED CLOSED BOMB DATA REDUCTION PROGRAM (CBRED)

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C. Price A. Juhasz

September 1977

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SCHEDULE 16. DISTRIBUTION STATEMENT (of this Report) Approved for public release; distribution unlimited. B. SUPPLEMENTARY 18. KEY WORDS (Continue on reverse side if necessary and identify by block number) Closed Bomb; Combustion; Burning Rates; Propellants - Burning Rates; Propellants - Closed Bomb Testing; Propellants - Combustion; Propellants -Artillery; Propellants - Small Arms. th. ABSTRACT (Combines on several shift If necessary and I-dentify by block number) (jmk) A versatile digital computer program was developed to provide linear burning rate information on propellants based on pressure-time records obtained from closed bomb firings. Some of the unique features of the program are: a treatment of heat loss based on radiative and convective heat transfer, capability of using single valued or tabular thermodynamic input, allowance for web deviation in the propellant sample, allowance for ignition deviation (flame spread) of the propellant sample, allowance for possible simultaneous burning of EDITION OF I NOV 60 IS QUICLETE UNCLASSIFIED

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propellant and ignition aid, allowance for tabular input of propellant surface versus mass fraction burned (to accommodate unusual geometries), and capability to treat vented vessel operations. The program was set up to operate on an interactive basis on a PDF 11/20 laboratory computer. In practice, once the program is called, the operator is guided in his choice of parameters (thermochemistry, heat loss, ignition deviation, etc.) by a series of prompts. A program overview is presented along with a description of equations, a derivation of the equations, and copies of program output and program listing.

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TABLE OF CONTENTS

	Page
ABSTRACT	1
	5
INTRODUCTION	11
OVERVIEW AND PROGRAM CAPABILITIES	14
THEORY	17
A. Simplified Derivation (1) Equation of State of Gas (2) Energy Balance Equation (3) Rate of Conversion of Solid to Gas (4) Linear Burning Rate	
B. Theory Used in the Program (1) Rate of Conversion of Solid to Gas (2) Linear Burning Rate (3) Heat Loss (a) Standard Option (b) Nonstandard Option	
PROGRAM STRUCTURE AND OPERATION	18
A. Standard Analysis	
B. Nonstandard Analysis	
C. Summary	
ACKNOWLEDGEMENTS	19
REFERENCES	10
APPENDIX A	S
APPENDIX B	9
APPENDIX C	9
APPENDIX D	1
APPENDIX E	9
	LIST OF SYMBOLS INTRODUCTION OVERVIEW AND PROGRAM CAPABILITIES THEORY A. Simplified Derivation (1) Equation of State of Gas (2) Energy Balance Equation (3) Rate of Conversion of Solid to Gas (4) Linear Burning Rate B. Theory Used in the Program (1) Rate of Conversion of Solid to Gas (2) Linear Burning Rate (3) Heat Loss (a) Standard Option (b) Nonstandard Option (b) Nonstandard Option PROGRAM STRUCTURE AND OPERATION A. Standard Analysis B. Nonstandard Analysis C. Summary ACKNOWLEDGEMENTS 3 APPENDIX A 4 APPENDIX C 5 APPENDIX C 5 APPENDIX D

																					Page
APPENDIX	F	•					•		•			•	•	•		•		•			83
APPENDIX	G	•	•	•	•	•	•	•	•		•		٠				•	•	٠	•	87
DISTRIBUT	ION	į į	LI	ST																	169

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Fortran Symbol	Symbol	Definition	Units
CAID	c _a	starting weight of ignition aid	1b
CPROP CNGWT	c _p	starting weight of propellant	1b
	g	gravitational constant (used as a conversion factor)	ft/sec ²
Н	h _c	convective heat transfer coefficient	Btu/in. ² -OR
SPACE2(3)	h _r	radiative heat transfer coefficient	Btu/in. ² -OR
TOTMOL	m _T	total moles of gas in chamber	lb-mol
RMOL	m _T	rate of change of total moles of gas in the chamber	mol/sec
RATE	r	linear burning rate of the propellant	in./sec
TIGN	t _{ig}	time of propellant ignition	msec
TMAX	t pa	time of maximum pressure	msec
Y(5)	w	weight of ignition aid burned	16.
Y(3)	wa1	weight of ignition aid combustion products in the chamber	1b
DY(5)	ů	ignition aid mass burning rate (w _a -AP ⁿ)	lb/sec
DY(3)	al	rate of change of ignition aid combustion products in the system (wal wa wa wa /ws)	lb/sec
Y(2)	w _i	weight of igniter combustion products in the chamber	1b

FORTRAN Symbol	Symbol	Definition	Units
DY (2)	wi	rate of change of igniter combustion products (w _i = - w _i w _n /w _s)	lb/sec
DNN	Wn	mass discharge rate through the nozzle	lb/sec
Y(6)	w _p	weight of propellant burned	1b
Y(4)	w _{pl}	weight of propellant combustion products in the chamber	16
DY (6)	w _p	mass burning rate of the propellant	lb/sec
DY (4)	w _{p1}	rate of change of propellant combustion products in the chamber (vented chamber operat (wpl = wp - wpl wn/ws)	lb/sec
Y(1)	w _r	weight of air in the chamber	1b
DY(1)	w _r	rate of change of air in the system $(w_r = -w_r w_n/w_s)$	lb/sec
FSYS	w	weight of gases in the chamber s=wr+wi +wal + wpl	16
DWSYS	w S	rate of change of the weight of gases in the system	lb/sec
DB	x	distance burned into the grain	in.
Y(8-12)	^x n	distance burned into the grain for the n-th charge increment (n has values of 1-5	in.)
SPACE1(10)	A _t	effective throat area(sonic control assumed)	in. ²
SPACE2(2)	A _w	bomb wall surface area	in. ²
СР	c _p	heat capacity at constant pressure	Btu/lb. OF

FORTRAN Symbol	Symbol	Definition	Units
CVP	c _v	heat capacity at constant volume	Btu/lb - °F
DCVP	ċ,	rate of change of heat capacity at constant volume	Btu/lb - OF-sec
CSTAR	C*	characteristic discharge velocity	ft/sec
OD	a	initial grain diameter	in.
PD	$\mathbf{p}_{\mathbf{p}}$	initial perforation diameter	in.
OOD	ים	instantaneous grain diameter (D' = D - 2x)	in.
PPD,PRFD	D'p	instantaneous perforation diameter (D ¹ _p = D _p +2x)	in.
E	E	end area of grain	in. ²
	E _e CA	energy of gases in the chamber	
	E*	energy from combustion of the propellant	
HTL1	HL	total heat loss to walls of the chamber	8tu
ARP	iL	instantaneous heat loss rate	Btu/sec
H	i.	average heat loss rate	Btu/sec
G	ii _{t.e}	convective heat loss rate	Btu/sec
HTL	ÅLT	radiative heat loss rate	Btu/sec
GL	L	initial grain length	in.
GGL `	L'	instantaneous grain length	in.

Variables used in derivation only

FORTRAN Symbol	Symbol	Definition	Units
XMA	Ma	molecular weight of ignition aid combustion products	NA
XMI	$M_{\mathbf{i}}$	molecular weight of igniter combustion products	NA
ХМР	M _p	molecular weight of propellant combustion products	NA
	$^{\rm M}_{ m r}$	molecular weight of air (defined as 29)	NA
XMW	M _s	molecular weight of gases in the system	NA
Y(7),P	P	pressure	psia
DP	ģ	rate of change of pressure with respect to time	psia/sec
PMAX	p m	experimentally measured maximum pressure in the system	psia
P	Pst	stagnation pressure	psia
РТНЕО	P _{Os}	theoretically computed maximum pressure in the system	psia
RSYS	Rs	gas constant for the system (R _s =R _u m _T /w _s)	in1b/1b ⁰ F
DR	Ř s	rate of change of gas constant for the system (R _s =R _u /w _s (m _T - m _T w _s /w _s)	in1b/1b- ⁰ F-sec
RU	Ru	universal gas constant	in1b/mo1 ⁰ F
AAB,S	s _t	instantaneous propellant surface area	in. ²

FORTRAN Symbol	Symbol	Definition	Units
AB	S _{tn}	instantaneous surface area of nth charge increment	in. ²
TSYS	T _s	gas temperature in the chamber (computed as ^O K; note also that ^O R= ^O F +459.69)	o _R
TSYS	. T _{st}	stagnation temperature	o _F
SPACE1(36)	Tw	bomb wall surface temperature	° _R
TAID, TS	T _{Oa}	isochoric adiabatic flame temperature of the ignition aid	o _R
T6,TZD	T _{Op}	isochoric adiabatic flame temperature of the propellant	° _R
BVOL	$v_{\mathbf{b}}$	empty bomb volume	in. ³
VSYS	V _s	system volume	in. ³
DVSYS	v _s	rate of change of system volume Vs=-wsn-nws+(wa+wp)/p	in. ³ /sec
WID, WOD	W	initial propellant web	in.
ALPHA	a	angle used in multiperforated grain surface calculations (MPGSC) defined in Appendix C	-
ВЕТА	3	angle used in MMGSC, defined in Appendix C	deg
GAN1	r	ratio of specific heats	NA
ETA	η	propellant covolume	in. ³ /1b
DETA	ή	rate of change of propellant covolume	in.3/1b-sec
ТНЕТА	9	angle used in MPGSC, defined in Appendix C	deg

Fortran Symbol	\$ymbo1	Definition	Units	
RHO	ρ	solid propellant density	lb/in. ³	
CPGM2	r	a function of the specific heat ratio defined in Appendix D	NA	

I. INTRODUCTION

Among the parameters of interest to the interior ballistician are the burning rate and chemical energy of the propellant used in propelling charges. These parameters must be known to predict the performance to be expected from a given gun-ammunition system. 1,2 Reliable data on the chemical energy of the propellant may be obtained from thermodynamic calculations based on the chemical composition and heats of formation of the propellant ingredients. 3,4 Propellant burning rates, however, cannot be predicted with similar reliability although various efforts at predicting burning rates from chemical compositions, 5,6,7 thermochemistry 8 and chemical kinetics have been and are being

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¹J. Corner, M. A., PhD, Theory of the Interior Ballistics of Guns, New York, John Wiley and Sons, Inc., 1950.

²Paul G. Baer and Jerome M. Frankle, "The Simulation of Interior Ballistic Performance of Guns by Digital Computer Program," Ballistic Research Laboratories Report No. 1183, December 1962, AD #299980.

A. C. Haukland and W. M. Burnett, "Sensitivity of Interior Ballistic Performance to Propellant Thermochemical Parameters," <u>Proceedings of the Tri-Service Gun Propellant Symposium</u>, 11-13 October 1972, Picatinny Arsenal, Dover, NJ, Vol 1, Section 7.3.

⁴Paul G. Baer, Ingo W. May, and Jerome M. Frankle, "A Comparison of Several Predictive Approaches in Charge Establishment for Large Caliber Artillery Systems," <u>Proceedings of the 11th JANNAF Combustion Meeting</u>, Jet Propulsion Laboratory, Pasadena, CA., September 1974, Vol 1; C.P.I.A. Publication 261, December 1974, The Chemical Propulsion Information Agency, Silver Spring, Md.

⁵William H. Tschappat, Lt.-Col, Ord. Dept., <u>Text-Book of Ordnance and</u> Gunnery, 1st Ed., New York, John Wiley and Sons, Inc., 1917, pp. 111-118.

⁶Albert A. Bennett, PhD, "Tables of Interior Bailistics," Ordnance Department Pamphlet No. 2039, April 1921.

P. W. Riefler and D. J. Lowery, "Linear Burn Rates of Ball Propellants Based on Closed Bomb Firings," Ballistic Research Laboratories Contractor Rerort No. 172, August 1974. (AD #921704L)

⁸Henri Muracur, "Chemie Physique sur la Reaction Entre La Temperature de la Explosion d' une Poudre et sa Viterse de Combostion," Comp. rend., Vol 187, 1928, pp. 289-294.

pursued. 9-15 (See, footnote,*). Useful burning rate data, therefore, must still be obtained experimentally.

Classically, two principal methods have been used to obtain burning rate data for propellants. These are the strand burner and the closed bomb. Both involve the burning of propellant samples in steel containers of sufficient strength to withstand high pressures. In a strand burner, a single sample (strand) of propellant is burned cigarette fashion under essentially a constant pressure. In practice, this is achieved by connecting a ballast volume to the combustion chamber so that the gases given off by the propellant contribute negligibly to overall system pressure. The regression of the burning surface is measured by timing the intervals between breaking of fuse wires embedded along the length of the propellant sample. Thus each experiment yields a single average value of linear burning rate for a given pressure. To obtain burning rates for a propellant over a range of pressures, a number of experiments are required. The resultant data are fitted to some mathematical form which then allows computation of the values of the burning rate at intermediate points. The method is straight-forward (though time consuming) and has been used for many years, especially by the rocket community.

^{*}Current Army efforts on . The stion modeling are centered in the Fundamentals of Combustic: Task of the Energetic Materials Technology Program of the US Army Materiel Development and Readiness Command, Alexandria, VA 22333.

 $^{^{9}}$ Ref 1, pp. 42-84 and reference therein.

¹⁰ Bryce L. Crawford, Jr., Clayton Huggett, and J. J. MoBrady, "The Mechanism of Burning of Double Base Propellants," <u>J. Phys.</u> and <u>Colloid Chem.</u>, Vol 54, 1950, pp. 854-862.

¹¹Robert G. Parr and Bryce L. Crawford, Jr., "A Physical Theory of Burning of Double Base Rocket Propellants," J. Phys. and Colloid Chem., Vol 54, 1950, pp. 929-954.

^{120.} K. Rise and Robert Ginell, "The Theory of Burning of Double Base Rooket Powders," J. Phys. and Colloid Chem., Vol 54, 1950, pp. 885-917.

¹³ R. E. Wilfong, S. S. Penner, and F. Daniels, "An Hypothesis for Propellant Burning," J. Phys. and Colloid Chem., Vol 54, 1950, pp. 863-872.

¹⁴ R. D. Geckler, "Mechanism of . mbustion of Solid Propellants," Selected Combustion Problems, London, Butterworths Scientific Publications, 1954, pp. 289-339.

¹⁵D. B. Spalding, "The Theory of Solid and Liquid Propellants," Combustion and Flame, Vol 4, 1960, pp. 59-76.

The second method of obtaining propellant burning rate data involves the closed bomb. In the closed bomb, a statistically adequate number of propellant grains are ignited and allowed to burn in a fixed volume under the pressure exerted by the propellant combustion gases. The pressure in the chamber builds up rapidly and is recorded as a function of time. From the pressure-time data it is possible to derive linear burning rate information for the propellant over a range of pressures from a single experiment, a marked advantage over the repeated tests required with the strand burner.* But where the closed bomb technique gains in experimental economy, it loses in the complexity of the data reduction method. The problem has been attacked in a variety of ways by a number of authors. 17-20 The earlier papers were aimed at providing methods for computing the data by hand. This lead to the use of a variety of simplifying assumptions both in the development of the theory and the form functions used. The more recent papers were aimed at computer solutions to the problem and, in general, provide a more complete treatment of the phenomenon. A brief bibliography of closed bomb burn rate reduction methods is included at the end of this report.

*Alternately, the data recorded may be the first derivative of pressure with respect to time vs. pressure. This is generally reduced to an average value of dp/dt (obtained at 0.250,0.3750,0.500, and 0.625 of the maximum pressure) which when compared with the value for a reference propellant (obtained under identical conditions) gives an idea of the burning characteristic to be expected of the sample. The "quickness" and "relative quickness" values so obtained can be quite useful for correlations with weapon performance characteristics.

¹⁶ Methoa 801.1.1, "Quiokness and Force Measurement of Propellant (Closed Bomb Method)," (Revised 21 Oct 75), Military Standard 286B, Department of Defense, Washington, DC 20301.

¹⁷C. M. Dickey, "Petermination of Burning Characteristics of Propellant," E.I. duPont de Nemours and Company, Explosives Department, Burnside Laboratory, Memorandum Report No. 31 (File BL-135-101), March 1943.

¹⁸ James H. Wiegand, "A Method of Calculation of the Burning Rate of Powders from dP/dt vs. P Records for closed Chambers," Ballistic Research Laboratories Report No. 546, June 1945.

¹⁹ E. Haeuseler and W. Dehl, "State of Development of Testing Procedures for Propellants in the Closed Vessel," Explosive toffe, Vol 18, 1970, pp. 41-52.

²⁰H. Krier and S.A. Shimpis, "Predicting Uniform Gun Interior Ballistics: Part I - An Analysis of Closed-Bomb Testing," University of Illinois Technical Report AAS 74-5, July 1874. Contract DAAD-21-73-C-0549.

The objective of the present work was to generate a comprehensive data reduction method providing more versatility than is available from any of the previous methods. The equations which were developed provide for the presence of igniter and ignition aid, as well as the ambient air present in the bomb. A more sophisticated heat loss treatment is included. Other unique features include allowance for changes in the thermodynamic characteristics of the combustion products and of web deviations in the propellant sample. In addition, the treatment allows for analysis of data from vented chamber experiments. The theory was implemented in a program written for an available laboratory computer. An attractive feature of the program is its user oriented "question and answer" format which allows the user to readily modify input variables and to make decisions on the various data reduction options. The reduced data are available for examination in easily interpretable format in a matter of minutes.

II. OVERVIEW AND PROGRAM CAPABILITIES

The program cupabilities are outlined in the following section. In the form listed here it requires as input a suitable data file of pressure vs time and first derivative of pressure with respect to time and a "header" or set of parameters describing propellant geometry and physical characteristics, propellant thermochemistry and experimental parameters.*

A listing of the input requirements is contained in Appendix A, program operation is interactive and is controlled from a terminal keyboard. Two modes of operation are possible, standard and nonstandard. The standard option provides a "normal" burning rate analysis on a more or less routine basis. This analysis assumes full burning of all of the ignition aid before ignition of the propollant, simultaneous ignition of all the propellant grains, a fixed set of dimen ions for all propellant grains, constant thermochemistry for the combustion products, and linear heat loss during the combustion of the propellant sample.

The second, nonstandard, mode of operation allows for operation of the program with a variety of options for special applications. A summary comparison of the "standard" and "nonstandard" modes of analysis is presented in Table I.

^{*} The data file is obtained by operating in the range data with a smoothing and differentiation program "SCHECK." The program and operations are to be described in a future report.

Table I. Comparison of Standard and Nonstandard Data Reduction Modes

是一个时间,我们就是一个时间,我们就是一个时间,我们就是一个时间,我们就是一个时间,我们也不是一个时间,我们也会会会会会会会会会会会会会会会会会会会会会会会会会 一个时间,我们就是一个时间,我们就是一个时间,我们就是一个时间,我们就是一个时间,我们就是一个时间,我们就是一个时间,我们就是一个时间,我们就是一个时间,我们就

NONSTANDARD MODE	Compute heat loss based on convective and radiative heat transfer coefficients.	Thermodynamic data input as table (function of pressure).	Statistical treatment of web deviations used in calculations	Allows definition of ignition deviation in terms of time to account for flame spread effects.	Propellant ignited before all of ignition aid is consumed.	Allows computation of mass discharge when pressure exceeds diaphragm burst pressure.	Allows input of any surface area function in tabular form.
OPTION NO.	1	2	ю	4	v	ý	7
STANDARD MODE	Constant heat loss rate throughout burning	Average values of impetus, covolume, molecular weight, and γ used.	Fixed web used in calculations	Simultaneous ignition of whole propellant charge.	All ignition aid burned before propellant ignited.	No provision.	Capabilities provided: sphere; cylinder; and single, seven, and nineteen- perforated cylinders.
FACILITY	Heat Loss	Thermochemistry	Propellant Geometry	Ignition	Ignition Aid	Vented Vessel	Burning Surface

Use of the program in either the standard or nonstandard mode has been simplified by allowing the use of propellant, igniter, and experimental data all from the same input file. Data on propellant geometry, thermochemistry, and dimensions are all recorded as "header" information at the time that the data are taken. (This is achieved using a separate program, CBDAP, closed bomb data acquisition program. The program will be described in another report.) The result is that conversion of the pressure-time data to linear burning rates may be performed simply without detailed knowledge of program operations. In performing a standard analysis, once the operator has entered the input data file, he is freed from providing any additional input except for specifying whether only the central portion (from 10 to 80 percent of maximum pressure) or all of the curve is to be plotted.

If changes are to be made in some of the propellant or igniter data or if special handling of the data is required, the nonstandard mode of analysis is employed. In this mode, the pertinent propellant and igniter data in the data file are displayed and the opportunity for changing or accepting the respective value is presented. In addition, the opportunity of selecting each of the options in Table I is presented to the user. Once these decisions have been made, data reduction proceeds as before.

Program output consists of: (a) a summary sheet or header describing the sample, experimental parameters and data on the maximum pressure and selected values of the derivative of pressure with respect to time (dP/dt), (b) a tabular listing of pressure, time, dP/dt, burning rate, web and surface fraction data, (c) superimposed plots of pressure (P) versus time (t) and dP/dt versus t. (d) a plot of dP/dt versus reduced pressure (P/P_{max}) , and (e) a log-log plot of burning rate as a function of pressure. The burning rate versus pressure plot includes a printout of the coefficient and the exponent in the equation

$$\mathbf{r} = AP^{n} \quad , \tag{1}$$

where: r = linear burning rate,

P = pressure,

A = burning rate coefficient, and

n = burning rate exponent.

obtained by a least squares fit of the data over the desired pressure range as well as statistical data on the "goodness of fit." An example of program output is given in Appendix B.

III. THEORY

A. SIMPLIFIED DERIVATION

The basic objective of the analysis is to derive linear burning rate data for propellant samples from pressure-time histories of their burning process in a closed bomb. Simply, the event involves the conversion of a solid sample composed of a large number of grains of a given geometry and size to a gas having a given amount of energy. Since the vessel is closed, the products may not escape, the pressure builds up and the propellant sample burns in an environment of a continuously changing pressure.

A variety of factors influence the conversion rate of the solid sample to gas. Those of primary interest are the propellant surface area, the pressure, and the propellant chemical composition. For all propellants, the conversion rate of solids to gas (i.e., the mass burning rate) is directly proportional to surface. For all gun propellants (and many others) the rate of regression of the propellant surface (the so-called linear burning rate, r) is directly proportional to pressure. As a general rule, the linear burning rates of propellants at given pressures are a function of the energy content of the composition.

To describe the process it is necessary to describe the gas produced, the unburned propellant and the energy balance for the process as a function of time. The derivation which follows is limited to a single propellant of constant thermochemistry. Eliminating complicating factors such as the presence of the igniter, ignition aid, etc., allows the generation of a simple instructive set of equations demonstrating the logic used. The actual equations used in the program are discussed in a later section and their derivation is given in the appendix.

(1) Equation of State of Gas

The combustion products may be described using the following equation of state:

$$PV_{s} = w_{p}R_{s}T_{s}$$
 (2)

where:

V_s = system volume

w_n = weight of propellant burned

 R_s = gas constant for the system

 $T_s = gas$ temperature in the chamber

The equation is formally analogous to the familiar Ideal Gas Equation. The difference is in the definition of the system volume term (V_s) which is defined as the chamber volume modified to reflect the presence of unburned propellant and the covolume correction, Equation (3)

$$V_{s} = V_{b} - \frac{c_{p}}{\rho} + \frac{w_{p}}{\rho} - w_{p}\eta$$
 (3)

where:

V_h = chamber volume

c_n = initial weight of propellant

 ρ = solid propellant density

η = propellant covolume.

It should be noted that the mixture of gases making up the propellant combustion products is treated as if it were a single gaseous species having specific properties of molecular weight, heat capacity and covolume. These properties, of course, are determined by the nature and stoichiometry of the combustion products.

Once the appropriate substitutions are made, the Equation of State becomes:

$$P\left[V_{b}-\frac{c_{p}}{\rho}+w_{p}\left(\frac{1}{\rho}-\eta\right)\right]=w_{p}R_{s}T_{s}$$
(4)

(2) Energy Balance Equation

The energy from combustion of the propellant sample is partitioned between the internal energy of the product gases and heat loss to the chamber. The Energy Balance Equation may be written as:

$$E_{cv} = E_{p} - H_{L} , \qquad (5)$$

where: E_{CV} = energy of gases in the chamber

 E_{p} = energy from combustion of the propellant, and

H_{I.} = heat loss to walls of the chamber

The equation may be rewritten as:

$$C_{\mathbf{V}^{\mathbf{W}}\mathbf{p}^{\mathbf{T}}\mathbf{s}} = C_{\mathbf{V}^{\mathbf{W}}\mathbf{p}^{\mathbf{T}}\mathbf{0p}} - \mathbf{H}_{\mathbf{L}}$$
 (6)

where: $C_V = \text{heat capacity at constant volume and}$

 T_{0p} = isochoric flame temperature of propellant.

$$H_{L} = \frac{C_{V}V_{s}}{R_{s}} (P_{0s} - P_{m})$$

 P_{0s} = Theoretically computed maximum pressure

 P_{m} = Experimentally measured maximum pressure

The temperature of the combustion gases, T_s , is less than the isochoric adiabatic flame temperature computed for the propellant composition, T_{Op} , due to heat losses to the walls of the chamber. The heat capacity at constant volume, C_V , is an average property between T_s and T_{Op} for the mixture of gases making up the combustion products of the formulation. It is defined per unit weight, rather than per mole.

(3) Rate of Conversion of Solid to Gas.

To obtain the equation for the rate of conversion of the solid propellant to gaseous combustion products, Equations (4) and (6) are differentiated. Differentiation of Equation (4) holding V_b , c_p , ρ , η and R_c constant, yields:

$$\begin{bmatrix} V_b - c_p + w_p & (\frac{1}{\rho} - \eta) \end{bmatrix} \frac{dP}{dt} + P & (\frac{1}{\rho} - \eta) & \frac{dw_p}{dt} = \\ R_s w_p & \frac{dT_s}{dt} + R_s T_s & \frac{dw_p}{dt} \end{bmatrix}$$
(7)

The rate of conversion of solid to gas is given by dwp/dt, the rate of formation of propellant combustion products. This is the term we are seeking to evaluate in terms of experimental parameters. To do this, it is necessary to define the rate of change of system temperature

(dT /dt). This is done by differenting the Energy Equation. Differentiation of Equation (6) holding C_{V} and T_{0D} constant and rearranging, yields:

$$\frac{dT_s}{dt} = \frac{(T_{0p} - T_s)}{w_p} \frac{dw_p}{dt} - \frac{\dot{H}_L}{C_V w_p}$$
 (8)

at this point, the following relationship is introduced:

$$R_s = C_V (\gamma - 1)$$

where:

y = ratio of heat capacities

The relationship is strictly true for ideal gases but is commonly used in describing real systems. It allows recasting Equation (8) in the following form:

$$\frac{dT_s}{dt} = \frac{(T_{0p} - T_s)}{w_p} \frac{dw_p}{dt} - \frac{\dot{H}_L}{R_s w_p}$$
(9)

Substituting the right hand side of Equation (9) into the differentiated Equation of State, Equation (7); yields:

$$\begin{bmatrix} V_b - \frac{c_p}{\rho} + w_p \frac{(1-\eta)}{\rho} \end{bmatrix} \frac{dP}{dt} = \begin{bmatrix} R_s T_s - P \frac{(1-\eta)}{\rho} \end{bmatrix} \frac{dw_p}{dt}$$

$$+ R_s w_p \begin{bmatrix} \frac{(T_{0p} - T_s)}{w_p} & \frac{dw_p}{dt} & -\frac{\dot{H}_L}{R_s w_p} \end{bmatrix}$$
(10)

The equation may now be solved for dw_p/dt , giving the rate of formation of propellant combustion products in terms of experimental data (P, dP/dt) and a series of constants (V_b , c_p , ρ , η , T_{0p} , R_s , H_L

and γ). The resulting equation is:

$$\frac{dw_{p}}{dt} = \frac{\left[V_{b} - \frac{c_{p}}{\rho} + w_{p} \frac{(1 - \eta)}{\rho}\right] \frac{dP}{dt} + \mathring{H}_{L} (\gamma - 1)}{R_{s}T_{0p} - P \frac{(1 - \eta)}{\rho}}$$
(11)

(4) Linear Burning Rate.

The linear burning rate, r, is the rate of regression of the propellant surface (dx/dt). To compute it one begins with considering the volume element burned through during an infinitesimal time interval. The following equation applies:

$$\frac{dw_{p}}{dt} = \rho S_{t} \frac{dx}{dt}$$
 (12)

where:

 S_{+} = the surface area of the propellant at any time t.

 $\frac{dx}{dt}$ = rate of regression of the propellant surface, equals r, $\frac{dx}{dt}$ the linear burning rate.

Equation (12) defines the rate of formation of propellant combustion products in geometric terms. If the propellant is composed of a number of identical grains, the propellant surface area may be computed using a variety of "form function" equations. Equations have been developed for spherical, cylindrical and perforated cylindrical (1, 7, and 19 perforations) grain geometries. The generalized equation is:

$$S_{t} = f(x) \tag{13}$$

where:

x = the distance burned into the grain.

When x equals zero, the surface area is the initial surface area of the propellant grain. As x increases positively, the surface area of the grain changes characteristically for each grain type. Form functions for all of the grain types mentioned above have been included in Appendix C.

Examining Equation (12), it is evident that the objective of computing the linear burning rate of the propellant from the closed bomb pressure-time data has been attained. The term dw_p/dt is defined by Equation (11), and the surface area, S_t , is defined by the Form Function Equation, Equation (13). This completes the simplified derivation intended for inclusion in the text.

B. Theory Used in The Program.

Both the event described and the equations describing it are considerably more complex than just described. In every experiment an igniter is used to start the propellant burning and in many an ignition aid is used as well. Further, the volume in the bomb not occupied by the propellant at the start of the experiment is occupied by ambient air. It is, therefore, evident that one is not dealing with the single component system described earlier. The Equation of State used in the program includes all components. The Energy Equation is also more complex, since the thermodynamics of the combustion gases are allowed to change and, in addition, allowance is made not only for heat loss from the system but mass loss as well (vented chamber operation). The Mass Burning Rate Equation derived for the system reflects these complexities. See Equation (27), Appendix D. Since the treatment allows for deviations in the ignition of the propellant charge as well as web deviations, computing the surface area at any instant is also slightly more complicated. Finally, the instantaneous Heat Loss Term (H,) is evaluated in the program in one of two ways, either as an average value, constant throughout the burn, or as a variable defined by the radiative and convective heat loss elements. The objective of this section is to provide some explanatory comments on several equations used in the program.

(1) Rate of Conversion of Solid to Gas.

For the purpose of discussion the Mass Burning Rate Equation derived in Appendix D has been regrouped and the numerator divided into a series of terms A through E. This is the form in which it appears below.

$$\frac{dw}{dt} = \frac{A + B + C + D + E}{P(\eta - \frac{1}{\rho}) + T_{0p}R_{s} + R_{u}T_{s} (\frac{1}{M_{p}} - \frac{1}{M_{s}})}$$
(14)

where:

R_u = universal gas constant

 M_p = molecular weight of propellant combustion products

 M_s = molecular weight of gas in the system

$$A = V_{s} \frac{dP}{dt}$$

$$B = \dot{H}_{L} (\gamma-1)$$

$$C = \begin{bmatrix} \gamma R_s T_s + Pn & w_{p1} \\ & w_s \end{bmatrix} \frac{dw_n}{dt}$$

where:

p₁ = weight of propellant combustion products in the chamber
(as opposed to the total weight of propellant burned)

 $\frac{dw_n}{dt}$ = gas discharge rate through the nozzle

$$D = w_s T_s \left[\frac{R_u}{(M_p)^2} \frac{dM_p}{dt} + \frac{R_s}{C_V} \frac{dC_V}{dt} \right] - Pw_s \frac{dn}{dt}$$

$$E = -\left[P(n-1) + T_{0a} R_{s} + R_{u} T_{s} \left(\frac{1}{N_{a}} - \frac{1}{N_{s}}\right)\right] \frac{d w_{a1}}{dt}$$

where:

 M_a = molecular weight of ignition aid combustion products

wal = weight of ignition aid combustion products in the chamber.

Several of the terms are associated with exercising program options previously described (Table I). The functions of Terms A through E are listed in Table II.

Table II. Functions of Terms A through E in Equation (14).

- A System volume term
- B Heat Loss.
- C Mass Loss.
- D Variable thermochemistry.
- E Contribution of simultaneously burning ignition aid.

The System Volume Term (A) is necessary for all computations. The Heat Loss Term (B) is included as required in the analysis. The term may be evaluated simply (standard option, Table 1) or comprehensively (nonstandard option). This will be discussed more fully in section B(3). In the case of vented chamber operation or in computing burn rates from artificially generated pressure time data, term B can be zero. The Mass Discharge Term, (C) is used in analyzing vented chamber experiments. Term (D), Variable Thermochemistry, is important in analyzing low pressure closed bomb data since the thermochemistry of the combustion products changes significantly with pressure at low pressures. The inclusion of the Ignition Aid Burning Term (E) becomes important when describing situations involving simultaneous burning of propellant and igniter. This is the case more often than not, though the decision on the overlap of ignition aid and propellant burning is made on the basis of experience by the program operator.

In essence, Equation (14) is Equation (11) appropriately modified to reflect the complexities of the experiment. This may be readily demonstrated by imposing the same assumptions on Equation (14) as were used in deriving the simplified Mass Burning Rate Equation (Equation 11). Assuming a single propellant (no igniter or ignition aid), constant thermochemistry, closed bomb operation (no mass loss through a nozzle) and the absence of air from the chamber, the following terms in Equation (14) may be eliminated:

- Term C. Under closed bomb conditions dwg/dt=0
- Term D. For constant thermochemistry dn/dt=0
- Term E. In the absence of an ignition dwal/dt=0 aid

$$\begin{bmatrix} R_u T_s & (\frac{1}{N_p} - \frac{1}{N_s}) \end{bmatrix}. \quad \text{for a single component} \\ \text{system} \qquad \qquad (1/N_p - 1/N_s) = 0$$

Elimination of the terms above reduces Equation (14) to Equation (11). The methods used in deriving the two equations were the same. The derivation of Equation (14) paralleling the approach used in the text for Equation (11) is given in Appendix D.

(2) Linear Burning Rate

The linear burning rate in the program is computed by Equation (12). Differences arise, however, in the computation of the instantaneous burning surface area S_{\downarrow} . Two of the program options treat ignition deviations in the propellant charge and web deviation in the propellant grains. Both options influence the burning surface area.

If an ignition deviation takes place, parts of the propellant charge begin to burn before others. In this case the simple computation of total burning surface area as a function of distance burned is not appropriate. What is done is to proportion the propellant charge into five parts (two each of 10 percent, two each of 20 percent and one of 40 percent) and to allow the ignition of the charge increments to differ by some arbitrar, time input by the operator. The distances burned into the surface of each portion of the charge are carried separately. Under these conditions the burning surface is computed according to the following equation:

$$b_{t} = \sum_{1}^{5} S_{t_{n}}$$
 (15)

where:

 S_{t_n} = surface area of the nth propellant charge increment

$$S_{t_n} = f(x_n)$$

x_n = distance burned into the surface of the nth propellant charge increment.

Of course, for simultaneous ignition of the propellant charge, Equation (15) reduces to Equation (13). It must be emphasized, however, that one has no a priori knowledge of the ignition deviation time of the propellant charge; so this treatment should be viewed with caution.

The web deviation is handled analogously; In this case, however, web deviation values may be obtained from actual measurements.

(3) Heat Loss

(a) Standard Option. Evaluation of the Heat Loss Term $(\mathring{H}_{_{\! 1}})$ is done in either of two ways. In the standard option it is some suitable average value, constant throughout the analysis. The following equation applies:

$$\dot{H}_{L} = \overline{\dot{H}}_{L} = \frac{C_{V}V_{s}}{R_{s}} \frac{(P_{0s}-P_{m})}{(t_{pm}-t_{ig})}$$
(16)

where:

H_L = average heat loss rate

Pos = theoretically computed maximum pressure (adiabatic conditions, contributions from propellant, igniter, ignition aid and air in system).

t_{nm} = time of maximum pressure

tig = time of ignition

The total heat loss is the difference between the adiabatically computed internal energy of the system and the internal energy computed from the maximum pressure observed. The total heat loss is converted into the average Heat Loss Rate ($\mathring{\text{H}}_{L}$) by dividing by the burning time interval ($t_{pm}^{-t}-t_{ig}^{-t}$).

(b) Nonstandard Option

In the nonstandard ortion, heat loss is analyzed into its convective and radiative components. It is assumed that during the time that the propellant burns the gas generation results in convective heat transfer to the chamber walls. This is, of course, accompanied by radiative heat losses as well. After the propellant is consumed it is assumed that the convective heat loss becomes insignificant relative to the radiative heat loss. These assumptions allow the following treatment.

(i) Radiative heat loss coefficient. The temperature of the gases in the system is given by:

$$T_s = \frac{P_s V_s}{R_s}$$

The radiative heat loss rate is given by:

$$\dot{H}_{Lr} = \frac{C_V V_S}{R_c} \frac{dP}{dt}$$
 (17)

where:

 \dot{H}_{Lr} = radiative heat loss rate,

Once a matched array of T_s , \mathring{H}_{Lr} data are generated, they may be fitted to the following relationship:

$$h_{\mathbf{r}}^{A}_{W} = \frac{\hat{H}_{L\mathbf{r}}}{4 \quad 4} \tag{18}$$

where:

h_r = radiative heat transfer coefficient

A = bomb surface area

 $T_{w} = bomb wall temperature.$

In carrying out these computations, T_w^4 may be neglected since it is indeed insignificant $(T_s^4>>T_w^4)$ in the radiant heat transfer. The bomb surface area (A_w^4) is one of the program inputs (alternately a default value of 18.1 in is available in the program) and, therefore, the evaluation of the Radiant Heat Transfer Coefficient (h_r^4) is accomplished. Throughout the analysis, h_r^4 is a constant.

(ii) Convective Heat Transfer Coefficient. The Convective Heat Transfer Coefficient (h_c) is assumed to be a function of the instantaneous mass flow as in the case of heat transfer in a pipe. The relationship is given as:

$$h_{c} = \frac{\left(dw_{S}\right)^{0.8}}{dt} \tag{19}$$

where:

h_c = convective heat transfer coefficient

 $F_c = proportionality constant$

The value of h is computed at every point in the analysis using the instantaneous mass generation rate. The value of the proportionality constant F is obtained using an approximation technique in which an approximate value of h is calculated from the average mass generation and heat loss rates and this approximate value is refined using the value of h and increments of $\Delta P/\Delta t$.

(iii) Heat Loss Rate, nonstandard option. To compute the Heat Loss Rate (\dot{H}_L) at any given point in the analysis the following equations are used:

$$\dot{H}_{LC} = h_C A_W (T_S - T_W) \tag{20}$$

where: \dot{H}_{LC} = convective heat loss rate

$$\dot{H}_{Lr} = h_r A_w (T_s^4 - T_w^4)$$
 (21)

$$\dot{\mathbf{H}}_{L} = \dot{\mathbf{H}}_{Lc} + \dot{\mathbf{H}}_{Lr} \tag{22}$$

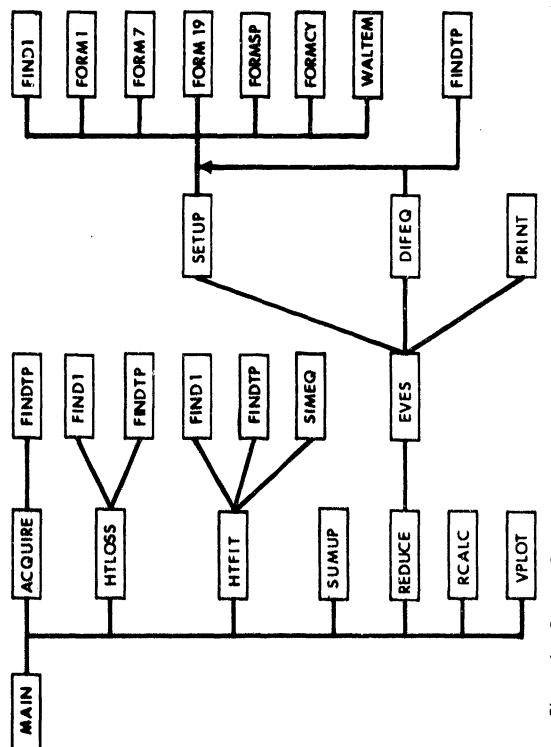
The initial value of the Wall Temperature (T_w) is assumed to be 450°K (this may be changed by the operator) and T_w is continuously adjusted throughout the analysis as is the Systems Temperature (T_s). The relationships governing heat conduction to the wall are given in Appendix E.

IV. PROGRAM STRUCTURE AND OPERATION

A. Standard Analysis

The program consists of a main program and a number of subroutines which handle reading in of data, unit conversions, heat loss computations, burning rate calculations, and printing and plotting of the output data. The program structure is outlined in Figure 1. The diagram indicates the relationship of the subroutines in the program. Capsule summaries of the functions of the subroutines are contained in Appendix F.

Program operation is most easily described by following a standard analysis sequence step by step. The options may then be seen as perturbations on the normal mode of analysis. A flow diagram of the program appears in Figure 2. The chart has been arranged to show the operations involved in a standard analysis in sequential form. The options possible, heat loss, mass loss, variable thermochemistry, etc., are shown offset from the main sequence of operations. The optional sections are marked with arterisks.



Interrelation of Subroutines. Program Structure (CBRED). Figure 1.

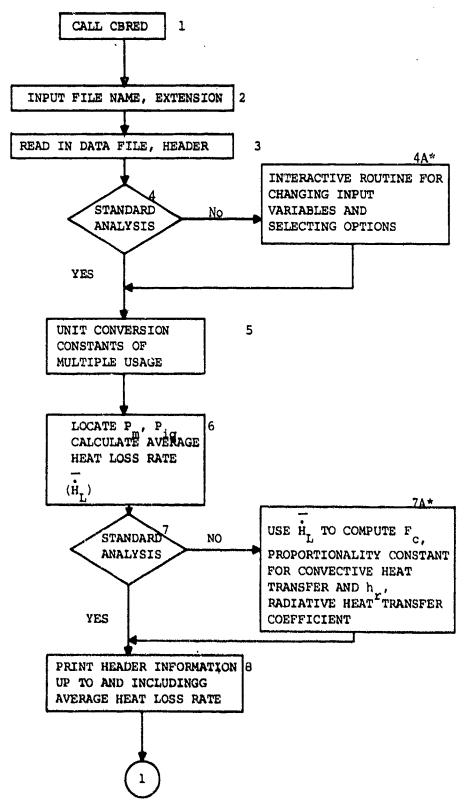


Figure 2-1. Closed Bomb Burn Rate Program (CBRED). Generalized Flow Scheme.

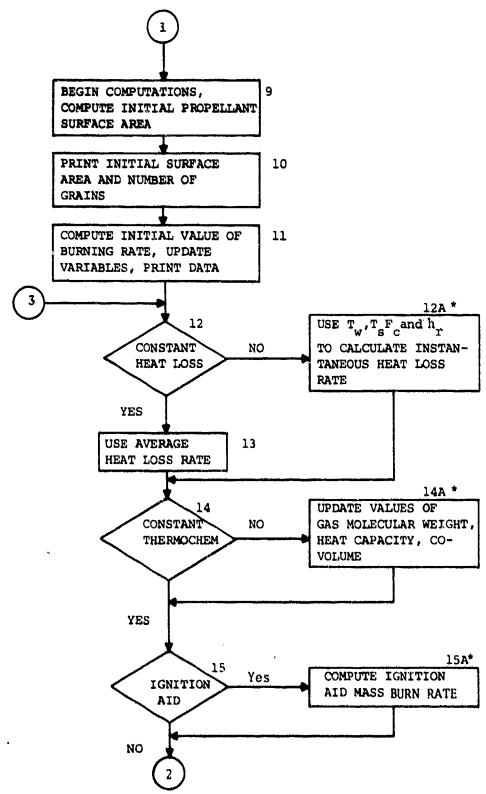


Figure 2-2. Closed Bomb Burn Rate Program (CBRED).
Generalized Flow Scheme

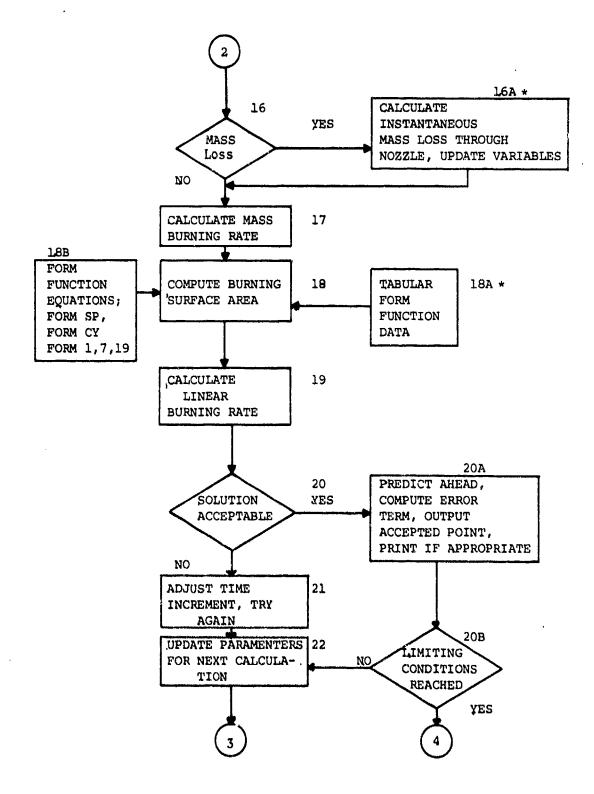


Figure 2-3. Closed Bomb Burn Rate Program (CBRED). Generalized Flow Scheme.

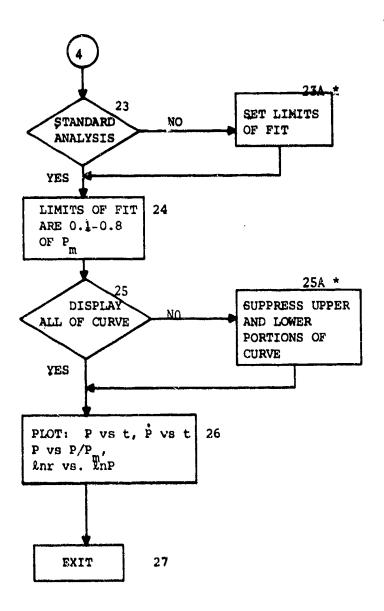


Figure 2-4. Closed Bomb Burn Rate Program (CBRED). Generalized Flow Scheme.

- 1. Start up (Blocks 1, 2 and 3). Once the program (CBRED) is called, it signs in and requests the file name and extension of the data file to be analyzed. On input of the file name and extension the data file is read in. The input data consists of matched arrays of pressuretime and time derivative of pressure-time data as well as a data set consisting of experimental, thermodynamic and grain geometry data (see Appendix A). The data is stored on magnetic tape and is handled as a single unit. (At the time the experiments run, the operator records in the data file all the parameters required for data reduction. This is done using a separate program, CBDAP. The pressure-time data obtained is differentiated by a second program, SCHECK. The "header" information is carried through to the new data file. Opportunities for updating the "header" information are provided by both programs). These operations take place in the main program and subroutine ACQUIRE.
- 2. Mode of Analysis (Block 4). The question is posed to the operator whether the analysis to be performed will be in the standard or nonstandard mode. Once the choice for the standard analysis mode has been made, the program goes into execution without further input from the operator. See subroutine ACQUIRE also.
- 3. Initialization (Block 5). At this point a variety of operations necessary to initialize the problem are performed. The input parameters are converted to a consistent set of units. At present a pound, foot (inch), second, BTU system is employed, but gas temperatures are in degrees Kelvin. A number of constants of multiple usage are also computed. This also takes place in subroutine ACQUIRE. On completion of this task, control reverts to MAIN which then passes operations to subroutine HTLOSS.
- 4. Heat loss (Block 6). At this point the input data file is searched for the maximum pressure, the maximum time derivative of pressure and selected values of P and dP/dt. The ignition pressure is computed using the igniter weight and thermochemistry and the theoretical (adiabatic) maximum pressure is calculated. This data is used to compute the Average Heat Loss Rate (H₁). Other operations including setting the ignition time for the analysis, setting the starting time interval, computing the initial system temperature and gas generation rate are also performed. These operations are performed in subroutine HTLOSS. Control then passes to MAIN which relinquishes operation to subroutine SUMUP.
- 5. Run header (Block 8). The function of subroutine SUMUP is to provide the identifying information for the run. Information on the data file designation, propellant parameters, igniter data, bomb data as well as summary information on the pressures and the time derivatives of pressure obtained are printed. The first page and part of the second page of Appendix B are the output from this operation. The only portion of the header not printed at this point relates to the initial

propellant surface area and the number of propellant grains. Subroutine SUMUP passes control back to MAIN which then passes control to subroutine REDUCE.

6. Begin Computation (Blocks 9-11). Subroutine REDUCE controls the differential equation solver, subroutine EVES. It specifies the number of differential equations to be solved, the stopping time and the interval at which data are to be printed out. (See pages 3 and 4 of Appendix B). Control then passes from REDUCE to EVES until the stopping time (or some other limiting condition) is reached.

To initiate computations, subroutine EVES calls subroutine SETUP whose job it is to provide the initial values of the functions and the factors needed to calculate the starting values of the differentials. Values are assigned to the propellant mass generation rate, ignition aid burning rate, mass loss rate, thermodynamic parameters, heat loss, etc. In a standard analysis, the initial heat loss rate is the average heat loss rate computed earlier in subroutine HTLOSS. The initial mass loss rate and igniter burning rate are zero. The initial surface area of the propellant is computed by calling the appropriate form function subroutine. The surface area data and the number of propellant grains are then printed, completing the "header" section of the program output (Block 10). Control passes from SETUP through EVES to DIFEQ where the initial values of the differentials are computed. (The functions evaluated using DIFEQ and EVES are designated as Y(N) and their differentials as DY(N); see the FORTRAN column in the List of Symbols.) In a standard analysis, the initial mass burning rate is set equal to the ignition aid mass burning rate at the point of ignition. The other mass differentials (and their integrals) are, of course, zero.

Using the starting values of the function (Y_0) and its differential (\dot{Y}_0) EVES computes an estimated value for the function at the completion of the first time interval (\dot{Y}_1) . This value is fed back to DIFEQ where it is used to obtain a new value of the differential (\dot{Y}_1) . The difference between (\dot{Y}_0) and (\dot{Y}_1) is examined and compared with an error level built into the program. If the value passes, the value of the function (Y_1) and its differential (\dot{Y}_1) are accepted, the time interval is incremented (Block 9) and the process is repeated. The initial integration steps are purposely kept small to establish the initial table of differences required by the predictor-corrector technique. Subroutine EVES has the built in capability to adjust the size of the time step used as the analysis progresses so that under conditions where the predictor values are better than required by the difference critereon, the time steps are increased in size, providing a savings of computation time. Time steps are, of course, also automatically reduced by the program as required.

- 7. Compute Burning Rates (Blocks 12 through 22). The process just described is followed to obtain the required values of linear burning rates throughout the analysis. At each step where a new value of (dw_/ dt) is required, subroutine DIFEQ computes the values based on the updated values of the parameters in the Mass Burning Rate Equation. The value is examined in EVES and accepted or rejected as necessary. Blocks 12 through 22 are involved in the process; Of these 12 through 19 take place in subroutine DIFEQ and the appropriate form function subroutine. In a standard analysis the heat loss rate is constant and the average heat loss rate is used. Since propellant thermodynamic characteristics are constant, no updating of values is done. The same is true for both the ignition aid burning and the mass loss terms. If the solution of the differential equation (as judged in EVES) is acceptable at any point, values for burning rate are accepted along with those of w and dw /dt. At given intervals during the analysis the data are printed out (EVES calls subroutine PRINT), see pages 50 and 51 of Appendix B. Tests for limiting conditions (web, all burned, pressure = P_m , and time = T_m) are made throughout the analysis and once one of the limiting conditions is reached, EVES returns control to REDUCE which returns control to MAIN.
- 8. Fitting the burning rate data (Block 25). Once the reduction phase is completed, the burning rate data between 0.1-0.8 of P_m is fitted to an equation of the form $r=AP^n$. This is done in subroutine RCALC.
- 9. Plotting of data and results. (25,26). The question is posed to the operator whether the lower and upper portion of the burning rate curve are to be suppressed. Generally this is done, since the most meaningful data is between 0.1 & 0.8 of P . After this decision, P vs.t; dP/dt vs t; dP/dt vs P/P and $\ln v$ vs. $\ln P$ are prepared and the program exits. The complete output package from the program may be seen in Appendix B.

B. Nonstandard Analysis

If the nonstandard mode of analysis is chosen, an interactive display is activated in subroutine ACQUIRE (Block (4A) which allows temporary modification of the pertinent ballistic information appearing in the data file. In addition to these changes, decisions concerning a variety of options have to be made. A brief discussion of each of the options, follows.

1. Heat loss. (Blocks 7A, 12 and 12A). In the nonstandard mode, the first option concerns the assignment of heat loss. The standard option, (Block 6). was described earlier. The nonstandard option (Block 7A) is a much more sophisticated treatment which assigns values to the convective and radiative compenents of heat loss based on the decay portion (after $P_{\rm m}$) of the firing record itself. The theory was discussed earlier. In the reduction phase of the program

the heat loss rate (\mathring{H}_1) is computed at each point in subroutine DIFEQ. Both the convective and radiative heat loss rates are functions of the gas and wall temperatures. Subroutine WALTEM is, therefore, invoked to provide current values of wall temperature as are needed in the analysis. This is shown schematically by Block 12A in Figure 2.

- 2. Propellant Geometry (Blocks 18A & 18B). The purpose of this option is to approximate the effect of real propellant geometry by using a distribution of web values rather than nominal web values. computations take place in DIFEQ and the appropriate form function subroutine. A web deviation value may be input in the nonstandard mode resulting in the proportioning of the propellant charge into five parts: two each of 10 percent, two each of 20 percent, and one each of 40 percent of the total charge with webs differing from the nominal by plus or minus suitable factors times the web deviation derived from a normal population distribution curve. These five propellants, equivalent in weight to but not having the same initial area or the same overall form function as the total charge considered as a nominal case are treated as separate entities in the burning area portion of the reduction and the instantaneous areas summed for calculation of the linear burning rate. This option eliminates sharp discontinuities of surface area at certain points. (Computer generated pressure-time curves using this approach have resulted in much closer agreement with dP/dt data taken from real propellant geometries.) Only the burning area portion of the analysis is affected.
- 3. Ignition Deviation. This option permits the simulation of flame spreading effects that may occur in an experimental firing. A population distribution similar to the one above (propellant geometry section) is used. The propellant is assumed to be composed of five samples, each ignited at a different time. (It is the value of the time interval that is input from the keyboard.) The distance burned for each fraction is calculated separately. These values of x are used to compute the area of the burning surface of the charge during the analysis. A linear interpolation method is used to remove gross discontinuities in the burning area between the times of ignition of two adjacent samples and upon burnout of the first sample. Only the burning area portion of the program is affected by this option. Operations take place in subroutine DIFEQ which calls the appropriate form function subroutine as required.
- 4. Ignition Aid. (Block 15, 15A). This option is useful in describing situations in which not only an igniter but also an ignition aid is present. Since the ignition aid actually has a finite burn time relative to the propellant, the option allows modification of the "ignition pressure" to account for the partial burnout of the aid material at the time that propellant ignition takes place. During the early phase of the firing, therefore, both propellant and ignition aid will be contributing to the mass generation rate, hence to dP/dt. The aid contribution must be evaluated separately. At present this is

accomplished by external input or by analysis of the dP/dt portion of the record just prior to the chosen ignition point (pressure), translating this dP/dt into an equivalent mass burning rate of the aid and fitting it to a power law curve. Evaluation of the instantaneous value of the ignition aid burning rate is performed in subroutine DIFEQ. The function is integrated in EVES. The computation of the mass burning rate (Block 17) is, of course, affected (see Equation 14 term E).

- 5. Variable thermochemistry. (Blocks 14 and 14A). This option allows the input of the required thermochemical information for the propellant in the form of a table of values versus pressure. The table entry takes place in subroutine ACQUIRE. Values for the rate of change of covolume, heat capacity etc. are computed in subroutine DIFEQ and used in calculating the mass burning rate of the propellant (also in DIFEQ).
- 6. Vented Chamber Operation. (Blocks 16 and 16A). This option allows analysis of firings made with "leaking" or vented vessels. The effective vent area and the blow-out pressure of the vent are necessary inputs and their presence will automatically result in analysis beyond the time of P_{\max} and the computation of the mass-discharge term in the differential equations. Instantaneous temperatures as well as the effects of changing molecular weight are used in determining this term.
- 7. Burning Surface. (Block 18A). In this option, an arbitrary total burning surface area as a function of some characteristic burning dimension may be input to the program if none of the available form functions appear suitable. The input consists of a table of points, and the instantaneous area is interpolated linearly from the information furnished. For seldom used unique geometries, it avoids the nuisance of recompiling the program to include a special form function generator.

C. Summary

A versatile closed bomb data reduction program has been developed to compute linear burning rate of propellants from the pressure-time histories of their burning process. In Section II, the capability of the program as seen from the user's point of view was discussed, and a comparison of the standard and nonstandard modes of analysis was made. An introduction to the theoretical treatment was made in Section III and the relationship between the equations used in the program and the simplified derived equation was examined. In Section IV, program structure and operation were examined and a flow chart of the program as well as a diagram of the subroutine hierarachy was given. Finally, the impact of each of the options on program execution was discussed. The derivation of the mass burning equations (Appendix D) as well as a program listing (Appendix G) are provided.

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APPENDIX A

Input Requirements for CBRED

Input Requirements for CBRED*

- 1. Run ID. Provides input file designation for program.
- 2. Propellant Data.
 - a. Type
 - b. Weight (grams)
 - c. Density (grams/cc)
 - d. Grain Type
 - e. Length, OD, ID (inches)
 - f. Inner Web, Outer Web (inches)
 - g. Theoretical Impetus (ft-lb/lb)
 - h. Flame Temperature (°K)
 - i. Average Molecular Weight of Products (grams/mole)
 - j. Covolume (cubic inch per pound)
 - k. Gamma (ratio of specific heat)
- 3. Igniter Data
 - a. Type
 - b. Weight (grams)
 - c. Impetus (ft-1b/1b)
- 4. Experiment Data
 - a. Bomb Volume (cc)
 - b. Bomb Temperature (%)
- * The input information listed is actually what is required by the program being documented. Conversion to metric units is being planned.

- 5. Data Arrays (digital)
 - a. Pressure-Time* (Kpsi, milliseconds)
 - b. Time Derivative of Pressure-Time* (mega psi, milliseconds)

^{*}The input information listed is actually what is required by the program being documented. Conversion to metric units is being planned.

APPENDIX B

Sample Output, CBRED*

^{*}Program output is currently not in metric units. Conversion is being planned.

	FEM 218-5-829; HTX 65% ZMICROW, 18% 18MICRON	B. BBBB.							•
CB1V75.027 LOYA STUDY RUM SIX 31 JULY 1975 AAJ/REB	1000 F784-35-1 FEM 218-5-82 12.50159 1.63888 330.50838 LOT 18A THYRON LESSITH HTM	CY 6.82178, 8.25808, 8.888 8.92178, 8.82178, 319502.06408 2288.84351 19.94808 31.59308 1.27188 FLAKE, NON-GRAPHITED, NON-PERFORRATED		Dupont 700x 0.50858 362539.02000		ea.90688 380.58883 PT-3 1.64328		24.58988 22.42971 8.92221	14.86122 25.42654 33.55888 11.36322
RUN ID: RUN TITLE: DATE: GPERATOR:	PROPELLANT DATA: TYPE: (FICHT (GYS): DENSITY (GYCE): ANTIAL TEPERATURE (DEG K): COT:	GRAIN TYPE: LENGTH.OD.ID (IM): INNER LEB.OUTER LEB (IM): THECRETICAL IPPETUS (FT-LB/LB): FLAPE TEPPERATURE: RVERIGE KULECULAR LEIGHT OF PROD: CO-WOLUTE (CU IN/LB): CATER (RATIO OF SP HTS):	IGNITER DATA:	TYPE: LEIGHT (CHS): IMPETUS (FT-LBAB):	EQUIPMENT DATA	BOYB YEYP (DEG K): GAUGE TYPE: CALIBRATION FACTOR (PC/PSI):	RESULTS:	THEORETICAL MAX PRESS (KPSIA): (DSERVED MAX PRESS (KPSIA): IGNITER PRESSURE (KPSIA):	IGNITION TIPE INFORMATION: THE TO SOX PROX (MSEC): THE TO SOX PROX (MSEC): THE TO 185% PROX (MSEC): THE FOOM 18% TO 98% PROX (MSEC):

ATTON:
INFORT
OUTCORESS

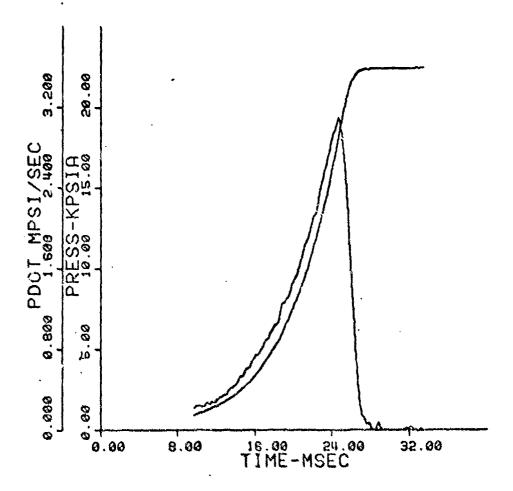
1.64.92 2.89854 2.54956 1.84129	3 89774 RECEDIED OF B - June 12
POOT AT .375 PPAX: PDOT AT .588 PPAX: PDOT AT .625 PPAX: AVERAGE PDOT:	MXIMM POST (POSI/SEC): 3

HEAT LOSS OPTION: CONSTRAT HEAT LOSS MURSER: 1453.51169

INITIAL SURFACE AREA (SO IN): 50.60207 KUREER OF GRAINS:

																																,																	
	FBEC-1	8.018	8.911	9.611	8.018	6.011	0.011	9.812	0.012	0.612	0.812	8.014	6.934	6.015	8.816	6.017	9.9.9	610.3	9.629	B.621	9.624	0.024	9.326	D.827	6.628	6.8 31	9.632	0.833	6.834	6.836	823.	9.079	9.641	0.643	9	0.e.	9.626	9.650	හ . පැත	9.057	6.83	9.062	9.063	9.868	0.871	9.04	2.0	. 462	0.083
	INCH BEEN	8.0000	6.6881	6.8881	. 0.6832	8.6863	6.6883	8.8884	6.6885	8.8986	0.000	6.8887	6.8898	6986.9	9.8918	0.8811	0.6512	6.00 13	0.8814	0.5815	0.6617	6.0018	6.0019	0.0621	9.0022	6.8824	9.8826	C. 6027	0.6629	6.6831	6.0833	0.6633	0.6837	6.6648	6.6842	6.8845	6.8847	0.000	6.8533	8.9556	6.6853	8.8962	5.8356	8.6869	8.6373	1.007		6.8083	6.0033 6.0033
	51.04 F.R	1.8000	6.9931	6.9981	8.3971	9.3962	9.9952	0.9941	6.9931	8.9928	6.3339	6.9838	9.9885	0.9873	9869	8.9846	6.9831	6.9916	8.9886	8.9784	8.9766	8.2747	8.9728	6.9767	8.9686	6.9663	9.9639	6.3513	92.88	6.9561	6 .9533	B.9584	0.9473	5.9442	9.948	8.9374	6.9339	5.9358	0.3259	8.3218	8.9175	6.9131	6.3834	928678	8.8386	6.8933	6.8878	6.8828	9.6768
1	ž	8.9559	6.8835	6.8873	6.8183	8.8146	6.6183	8.8224	6.6263	6.8384	0.2345	6. £388	6.8435	6.8482	6.6532	6.6562	6.6638	6.6695	6.8755	8.8316	6.8883	8.8334	8. 1826	6.1164	0.1184	8.1269	6.1339	8.1453	6.1548	B. 1647	8.1758	6. 1838	8.1978	6.2887	6,2289	0.2335	8.2464	e. 268%	0.2754	0.2594	6.3355	8.3218	8.3365	6.3559	6.27.38	B. 1703	6.4121	. 4.73	8.4537
	IN SEC	0.127	6.136	6.138	8.134	6.137	6.146	6.143	6.158	4.154	8.155	8.168	6.173	0.183	6.187	0.2B1	0.214	6.219	1.227	8.248	6.256	B. 256	96.7.98	8.323	6.313	B. 135	0.353	8.353	8.373	1.332	- 100 - 100	8.431	8.446	8 .466	8.4 89	6.561	6.538	. 596	6.535	6.612	9 .679	6.653	8.696	6.722	.7.5	.31		6.874	6.987
	PSI SEE	0.216	6 .236	0.241	6.233	6.237	0.257	0.262	8.267	6.273	6.278	6.389	6,325	9.342	0.351	8.382	6.413	9.6	6.443	8.472	.23	6.531	6.253	0.612	6.636	6.685	9.726	8.745	0.769	6.812	970.0	9.826	8.923	6.974	1.824		1.111	1.237	1.25	1.289	1.325	1.395	1.467	1.528	70%			1.833	1.833
1	PRESS KPS1A	6.922	6.978	1.638	1.633	1.156	1.215	1.284	1.348	1.417	.63:	1.256	1.637	1.719	1.838	1.895	1.998	2.181	2.211	2.322	2.467	2.534	2.717	2.678	3.823	3.191	3.365	3.551	3.738	1.937	4.143	4.363	4.589	4.838	5.677	5.339	5.665	5.693	6.213	6.1.3	6.835	7.193	7.551	7.925	9.311	1.717	9.14	565.6	16.639
101V73.00	11 12 12 13 14 14 14 14 14 14 14 14 14 14 14 14 14	9.736	9.986	18.236	10.486	10.736	16.986	11.236	11.486	11.736	11.500	12.236	12.486	12.736	12.996	13.236	13.486	13,736	13.986	14.236	14.466	14.735	14.986	15,236	15.486	15.736	13.986	16.236	16.486	16.735	986.91	17.236	17.486	17.736	17.966	10.236	18.486	18.736	10.935	19.235	19,436	19,736	19.985	28.236	22.456	22.736	38.82	21.236	21.486

	FEEC-1	6.689	8.695	8.832	69.698	9.184	9.169	. 113	6 .118	9.122	(C)	e.131	9 .135	9 .138	9.134	9 .126	6,113	6.835	9.874	9.626	6.637	9.621	0.011	8.886	6.684	6.683	9 .685	-8.533	-0.000	8.885	9.662	-6. B38	8.888	8.878 8	6.688	6.08G	6.888	6.666	6.686	6.886	888.8	6.891	6.681	B . 882	8.688	6.689	-8.888	9. 001
700	INCH INCH INCH	8.5894	6.6099	6.916.4	9.8183	6.6115	6.0128	2210.0	6.0133	8.8139	57.5.6	6.8154	0.0161	6.9169	9.8126	6.0184	6 .0191	8.8197	8.8282	8.8286	0.6283	8.8218	8.6212	8.8212	6.0213	8.8213	8.8214	2.6214	· 8.8214	0.8214	6.6215	6.8215	6.8215	6.5215	6.8216	9.62 is	8.6216	0.6216	6.8217	6.8217	8.8217	0.0217	6.6217	6.6218	8.6218	6.8218	0.8218	8.8219
8		6.8838	0.8632	a.8565	8.8496	6.8423	8346	9929	6.9133	6.8836	8.5358	6.7912	9.7916	6.7717	0.7618	6.7523	6.7416	6.7359	6.72% 6.72%	8.7246	8.7218	8.7187	6.7173	6.7165	6.7158	6.7153	6.7149	24.5.00	0.7142	6.7139	6.7134	6.7131	6.7128	e.7	8.7123	6.7128	6.7117	0.7114	0.7112	8.7189	6.7186	6 .6363	6. Bees	8.86 83	9.0da	6.6089	6 .8668	9.603.B
į	ž	6.4755	6.4382	6.5217	5456	C. 1.	6.5978	5.624	6.6522	6.6813	6.7114	0.7464	6.7.30	8.885Z	8.8383	6 .8687	0.000	6.9211	8. 948G	6.9863	9.9676	6.9747	9791	8.9817	6.9838	6.9853	8.9868	87.00°C	8.3887	8.9897	6.9912	6.9922	8.993	6.25	000.00 000.00	8.9956	90000 B	8.397A	6.9992	1666.3	6565.8	1.8818	1. 86 229	1.8831	1.6542	1.8851	1.6859	1.0253
1400	THE SET	8.945	6.992	1.673	. 8556 	1.153	5,183		1.293	1.349	1.416	1.463		1.557	1.521	1.446	e)	 	6.837	8.683	6.463	6.286	6.163	9.116	260.0	8.674	6.673	00.00 10.00	6.84J	8.878	8.878	6.643	6 .663	0.00	. e. c.	9.00	9. 9.	. 44 to 1	. e. d.	5.0 (3.0	6.847	6.947	6.047	6.647	6.04°	0.847	6.847	6.847
100	PSI/SEC	1.975	2.869	2.127	2.136	2.34	2.438	250.7	2.6	2.745	2.023	2.34	3.622	3.638	3.662	2.858		2.125	1.558	1.258	6.822	6.474	B.244	₽. T	8.693	8.6±8	8.856	100 d	-0:863	6.649	6.838	-6.663	3.84G	5 5 5	6.858	6.658	6.688	6 .	8.88.8 8.88	9.8£	6.683	9.854	6.614	8.634	6.657	6.894	-6.885 885	6.513
27 DECCEC	KP3IA	19.542	11.648	11.577	12.118	279.75	13.276	13.388	4.54	13.221	10. V	15.623	17.336	19.166	18.934	13.667	07 to 12 to	28.928	21.402	21.769	22.828	22.184	22.272	22.313	22.345	22.368	22,378	22.331	22.391	•	22.483	22.485	22.765				22.45	22.432	22.485	22.485	22.484	27.418	22.413	22.421	22.425	22.428	22.425	20.407
CB1775.827	111	21.736	2:.986	22.236	22.486	20.00	28.77	63.630	23.485	23.736	23.3.3	24.236	24.48 6	24.736	24.986	23.236	23.45	25.736	23.936	26.236	26.486	26.735	26.986	27.236	27.486	27.736	27.986	20.75	23.488	29.736	28.986		٠	23	23.996	٠	13.486	18.736	18.998	31.236	31.486	31.736	•	12.236	22.48 6	32.736	12.585	33.236

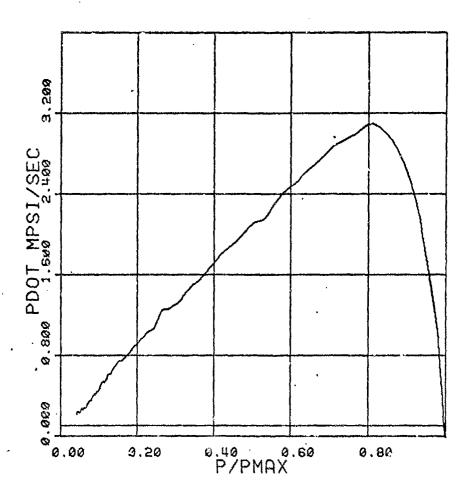


MUN ID: 051 V76.027

RLN TITLE: LOVA STUDY RUN BIX

PROP TYPE: LOWA F784-S6-1 FEN 216-6-029; HOX 66% SHICKON, 16% IGHICKON

ORGIN TYPE: OY



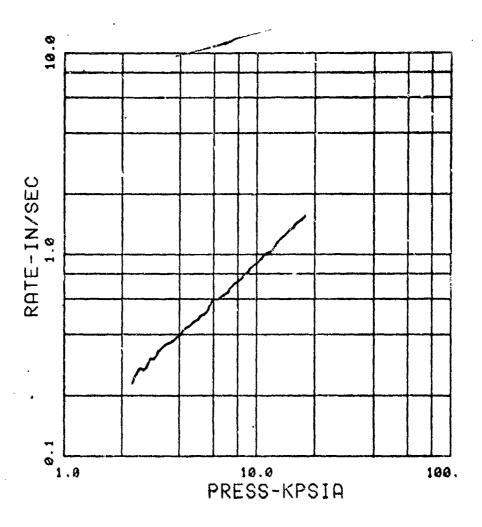
RUN ID: 081775.027

RUN TITLE: LOVA STUDY RUN BIX

PROP TYPE: LOVA F704-35-1 FEM 218-5-029; HAX 65% 2HICRON, 19% 10HICRON

GRAIN TYPE: OY

POUT	AT	PIPMAX
1.112		5.697
1.806		6.411
2.099		11.215
2.559		14.019



RUN ID: 031V75.027

RIN TITLE: LOVE STILDY RUN BIX

PROP TIME: LOVE F704-30-1 FEM 218-6-020; MIX 86K 2NICRON, 19K 10MICRON

GRAIN TYPE: CY

THE CONSTRUTE IN THE EQUATION R . ROPOON ARE

Ri 8.996233

1 **6.696**

FOR P/PHRX & 100 TO 0.800

COCFFICIENT OF DETERMINATION : 0.9999

MER CENT ROOT HEAN ERROR : 1.6034

APF IDIX C

Form Function Equations used in CBRED

Equation for FORM SP

Form Function of a Sphere. S, Surface Area, as a function of x, The Distance Burned

W = D

where: D = initial diameter of sphere

W = propellant web

x = depth burned at time t.

S = 0 for $W \le 2x$

Otherwise:

D' = D-2x

Surface Area

 $A=4\pi\left(\frac{D!}{2}\right)^2$

Equations for FORM CY

Form Function of a Right Circular Cylinder. S, Surface Area as a Function of \mathbf{x} , the Distance Burned

W = D

where: D = initial grain diameter.

W = propellant web

x = depth burned at time t

S = 0 for $L \le 2$ x (L = initial grain length)

S = 0 for $W \le 2$ x

Otherwise:

D' = D-2x L' = L-2x

End Area:

 $E = \pi/4 (D^*)^2$

Surface Area:

S = 2 E+m L' D'

Equations for FORM1

Form Function of a Single Perforated Right Circular Cylinder (Axially Symmetrical). S, Surface Area, as a function of x, the Distance Burned.

$$W = \frac{D - D_p}{2}$$

where:

And the second second second second second

W = propellant web

D = initial grain diameter $D_p = initial perforation diameter$

x = depth burned at time t.

S = 0 for L < 2x (L= initial grain length) S = 0 for W < 2x

Otherwise:

$$D^{\dagger} = D-2x$$

$$L^{\dagger} = L-2x$$

$$D_{p^{\dagger}} = D_{p} + 2x$$

End Area:

$$E = \frac{\pi}{4} [(D')^2 - (D_{p'})^2]$$

Surface Area:

$$S=2E + \pi L^{*} (D^{*} + D_{p^{*}})$$

Equations for FORM 7

Form Function of a Seven Perforated Right Circular Cylinder (Axially Symmetrical). S, Surface Area, as a Function of x, the distance burned.

I. To splintering:

$$W = D - 3 D_{p}$$

where:

W = propellant web
D = initial grain diameter
D_D = initial perforation diameter

x = depth burned at time t

S = 0 for $L^{1} \le 2x$ ($L^{1} = instantaneous grain length)$

 $S = S_{max}$ for W=2x

 $S = < S_{max}$ for W<2x

for 0 < x, define

$$\begin{array}{l} D^{+} = D-2x \\ L^{+} = L-2x \\ \qquad \qquad (L = initial \ grain \ length) \\ D_{\mathbf{p}^{+}} = D_{\mathbf{p}} + 2x \end{array}$$

Then, for $0 \le 2x \le W$

End Area

$$E = \frac{\pi}{4} [(D')^2 - 7(D_{p'})^2]$$

Surface Area:

$$S = 2E + \pi L' (D' + 7 D_{p'})$$

Form Function of a Seven Perforated Right Circular Cylinder (Axially Symmetrical). Surface Area, S, as a Function of x, The Distance Burned (continued).

II. After Splintering*

define: W_w=D_p+W and let

C=min
$$\left\{L, \frac{D^2 - D_p^2 + 4W_w^2 - 2W_w^D \sqrt{3}}{2(D + D_p - W_w \sqrt{3})}\right\}$$

Then for $W<2x\le C$, let

$$\Theta = 2 \cos^{-1} \left\{ \min \left(\frac{W_{w}}{D_{p'}}, 1 \right) \right\}$$

$$\alpha = \cos^{-1} \left\{ \min \left(\frac{1/4 \left[(D')^{2} - (D_{p'})^{2} \right] + W_{w}^{2}}{W_{w}^{D'}}, 1 \right) \right\}$$

$$\beta = \cos^{-1} \left\{ \max \left(\frac{1/4 \left[(D_{p'})^{2} - (D')^{2} \right] + W_{w}^{2}}{W_{w}^{D}}, -1 \right) \right\} - \frac{\Theta}{2} - \frac{\pi}{3}$$

 E_1 = End area of outer slivers for $\alpha < \pi/6$

$$E_1 = 3 D' W_w \sin \alpha + 3/2 \left[(D')^2 (\pi/6-\alpha) - W_w^2 \sqrt{3} - (D_p)^2 (\beta+1/2 \sin \theta) \right]$$

 $E_1 = 0$ for $a > \pi/6$

S₁ = Surface area of outer slivers

for a< 11/6

$$S_1 = 2E_1 + (6\beta Dp' + (\pi - 6\alpha) D') L'$$

$$S_1 = 0$$
 for $\alpha \ge \pi/6$

* Treatment developed by Mr. Franz Lynn, USABRL.

 E_2 = End area of inner slivers for $0 < \pi/3$

$$E_2 = 3/2 \left[w_w^2 \sqrt{3} - 3/2 (D_{p_1})^2 (\sin \theta + \pi/3 - \theta) \right]$$

 $E_2 = 0 \text{ for } \Theta \ge \pi/3$

 S_2 = Surface area of inner slivers for $\theta < \pi/3$

$$S_2 = 2E_2 + 9D_{p'} (\pi/3-\theta)L'$$

$$S_2 = 0 \text{ for } \pi/3 \leq 0$$

S = Total surface area of slivers

$$S = S_1 + S_2$$

E = Total end area

$$E = E_1 + E_2$$

for C < 2 x, E=0, S=0

Equations for FORM 19

Form Function for a Nineteen Perforated Right Circular Cylinder (Axially Symmetrical). S, Surface Area, as a Function of x, The Distance Burned.

To Splintering

$$W = \frac{D-5}{6} \frac{D_p}{P}$$

where:

W = propellant web D = initial grain diameter

 $D_{\mathbf{p}} = \mathbf{initial}$ perforation diameter

x = depth burned at time t

S = 0 for L' < 2x

(L' = instantaneous grain length)

S = S for W = 2x $S < S_{max}^{max}$ for W < 2x

for $0 \le x$, define:

D' = D-2x

 $L^{\dagger} = L-2x$

(L=initial grain length)

 $D_{p*} = D_p + 2x$

Then, for $0 \le 2x \le W$

End Area

$$E = \frac{\pi}{4} [(D')^2 - 19 (D_{p'})^2]$$

Surface Area

$$S = 2E + \pi L' (D' + 19 D_{p'})$$

For Function of a Nineteen Perforated Right Circular Cylinder (Axially Symmetrical). Surface Area S, as a function of x, The Distance Burned. (continued).

II. After Slivering*

define: $W_W = D_p + W$ and let

$$C = \min \left\{ L, 1/2 \left[D - D_p - W_w \sqrt{\frac{12[(D + D_p)^2 - 16W_w^2]}{[D + D_p]^2 - 12W_w^2}} \right] \right\}$$

Then, for w<2x<C, let

$$\Theta = 2 \cos^{-1} \left\{ \min \left(\frac{W_{w}}{D_{p}}, 1 \right) \right\}$$

$$\alpha_{1} = \cos^{-1} \left\{ \min \left(\frac{(1/8 [(D')^{2} - (D_{p'})^{2}] + 2 W_{w}^{2})}{W_{w}D'}, 1 \right) \right\}$$

$$\alpha_{2} = \cos^{-1} \left\{ \min \left(\frac{(1/4 [(D')^{2}] - (D_{p'})^{2} - (D_{p'})^{2}] + 3 W_{w}^{2}}{W_{w}D'}, 1 \right) \right\}$$

$$\beta_{1} = \cos^{-1} \left\{ \max \left(\frac{(1/8 [(D_{p'})^{2} - (D')^{2}] + 2 W_{w}^{2}}{W_{w}D_{p'}}, -1 \right) \right\}$$

$$\beta_{2} = \cos^{-1} \left\{ \max \left(\frac{(1/4 [(D_{p'})^{2} - (D')^{2}] + 3 W_{w}^{2}}{W_{w}D_{p'}}, -1 \right) \right\}$$

and:

$$\alpha = \alpha_1 + \alpha_2$$

 $\beta = \beta_1 + \beta_2 - \theta - 5\pi/6$

* Treatment developed by Mr. Franz Lynn, USABRL.

 E_1 = End area of Outer Slivers for $\alpha < \pi/6$

$$E_{1} = 3 D' W_{w}(2 \sin \alpha_{1} + \sqrt{3} \sin \alpha_{2}) - 6 W_{w}^{2} \sqrt{3} + 3/2 [(D')^{2} (\pi/6-\alpha) - (D_{p'})^{2} (\sin \theta + \beta)]$$

 $E_1 = 0$ for $\alpha > \pi/6$

 $S_1 = Surface Area of Outer Slivers$ for $\alpha < \epsilon \pi / 6$

$$S_1 = 2E_1 + 6 [D'(\pi/6-\alpha) + D_p' \beta] \cdot L'$$

 $S_1 = 0 \text{ for } \alpha \ge \pi/6$

 E_2 = End area of Inner Slivers for $0 < \pi/3$

$$E_2 = 6 \left[W_W^2 \sqrt{3} - 3/2 \left(D_{p_1}\right)^2 \left(\sin \theta + \pi/3 - \theta\right)\right]$$

 $E_2 = 0 \text{ for } \theta \ge \pi/3$

S₂ = Surface Area of Inner Slivers

$$S_2 = 2E_2 + 36 D_{p'}(\pi/3 - 0)L'$$

$$S_2 = 0$$
 for $0 > \pi/3$

S = Total Surface Area of Slivers $<math>S = S_1 + S_2$

E = Total End Area $E = E_1 + E_2$

for C<2x, E=0 and S=0

APPENDIX D

Derivation of Mass Burning Rate

Equation for CBRED

Derivation of the Mass Burning Rate Equation for CBRED

The same order of presentation is followed as in the text. The Equation of State is presented first, the Energy Equation, second, followed by the Mass Burning Rate Equation.

(1) Equation of State of Gas

$$PV_{s} = W_{s} R_{s} T_{s}$$
 (23)

where:

The state of the s

P = pressure

$$V_s = \text{system volume}$$

= $V_b - \frac{(c_a + c_p)}{\rho} + \frac{(w_a + w_p)}{\rho} - w_s \eta$

V_b = empty bomb volume
c_a = starting weight of ignition aid
c_p = starting weight of propellant
p = solid propellant density (assumed same for ignition aid)
w_a = weight of ignition aid burned*
w_p = weight of propellant burned*
w_s = weight of air in thamber
w_s = weight of initiator combustion products in chamber
w_a = weight of ignition aid combustion products in chamber*
w_a = weight of propellant combustion products in chamber*
n_a = covolume

$$R_s = \frac{R_u m_T}{w_s}$$

R₁₁ = universal gas constant

 M_r = molecular weight of air (taken as 29.)

M_i = molecular weight of initiator combustion products

 $M_{\rm a}$ = molecular weight of ignition aid combustion products

M_D = molecular weight of propellant combustion products

* Note: For a closed bomb system there is an obvious redundancy between W_p and W_{p_1} and W_a and W_{al} . But for a leaking, or vented vessel, the distinctions are important.

(2) Energy Balance Equation

It is more convenient to describe the energy balance dynamically. The following governing equation applies:

$$\frac{d(C_{V}w_{s}T_{s})}{dt} = C_{V} [T_{0a} \hat{w}_{a} + T_{0p} \hat{w}_{p}] - \hat{H}_{L} - C_{p} T_{s} \hat{w}_{n}$$
 (24)

where:

C_V = heat capacity at constant volume (assumed same for ignition aid)

 T_{0a} = isochoric adiabatic flame temperature of the ignition aid

 T_{0p} = isochoric adiabatic flame temperature of the propellant

 $\dot{\mathbf{w}}_{g} = AP^{n}$ By definition. Ignition aid mass burning rate.

 \dot{v}_{n} = mass burning rate of the propellant.

 \dot{H}_{t} = heat loss rate

 $C_{\rm p}$ = heat capacity at constant pressure

$$\dot{\mathbf{w}}_{n} = \mathbf{g} \frac{\mathbf{p}}{\mathbf{st}} \frac{\mathbf{A}_{t}}{\mathbf{C}^{*}}$$

where:

g = gravitational constant

P_{st} = stagnation pressure

 A_{+} = effective throat area (sonic control assumed)

C* = characteristic discharge velocity

$$= \left[\frac{g R_u T_{st}}{\tau^2 M_s} \right] 1/2$$

T_{st} = stagnation temperature

M_s = system molecular weight

 τ^2 = a function of the specific heat ratio

$$\left(\frac{\gamma+1}{\gamma-1}\right)$$

$$\tau^2 = \gamma\left(\frac{2}{\gamma+1}\right)$$

(3) Rate of Conversion of Solid to Gas

Solving the Equation of State (23) for T_s and differentiating yields:

$$\frac{dT_s}{dt} = \frac{1}{R_s w_s} \left[P\dot{V}_s + V_s \dot{P} - \frac{PV_s}{R_s w_s} (w_s \dot{R}_s + R_s \dot{w}_s) \right]$$
(25)

where:

$$\dot{V}_s = -w_s \dot{\eta} - \eta \dot{w}_s \left(\frac{\dot{w}_a + \dot{w}_p}{\rho} \right)$$

$$\dot{\eta} = \frac{d\eta}{dP} \times \frac{dP}{dt}$$
 By definition

$$\dot{R}_{s} = \frac{R_{u}}{w_{s}} \left[\dot{m}_{T} - \frac{m_{T} \dot{w}_{s}}{W_{s}} \right]$$

$$\dot{m}_{T} = \dot{w}_{r} + \dot{w}_{i} + \dot{w}_{a} + \dot{w}_{p} - \frac{w_{p} \dot{w}_{n}}{M_{p} w_{s}}$$

$$-\frac{\mathsf{W}_{\mathsf{p}1}\mathsf{M}_{\mathsf{p}}}{(\mathsf{M}_{\mathsf{p}})^2}$$

$$\dot{w}_{r} = -\frac{w_{r}}{W_{s}} \times \dot{w}_{n}$$
 By definition

$$\dot{w}_{i} = -\frac{w_{i}}{w_{s}} \times \dot{w}_{n}$$
 By definition

$$\dot{w}_{al} = \dot{w}_a - \frac{w_{al}}{w_s} \dot{w}_n$$
 By definition

$$\dot{\mathbf{w}}_{s} = \dot{\mathbf{w}}_{r} + \dot{\mathbf{w}}_{i} + \dot{\mathbf{w}}_{al} + \dot{\mathbf{w}}_{p} - \underbrace{\mathbf{w}_{pl}}_{\mathbf{w}_{s}} \dot{\mathbf{w}}_{n}$$

Note also:

$$\dot{\mathbf{w}}_{pl} = \dot{\mathbf{w}}_{p} - \frac{\mathbf{w}_{pl}}{\mathbf{w}_{s}} \dot{\mathbf{w}}_{n}$$
 By definition

Differentiating the left hand side of Equation (24) as indicated gives:

$$C_{V} \overset{\bullet}{w_{s}} \overset{\bullet}{T_{s}} + C_{V} \overset{\bullet}{T_{s}} \overset{\bullet}{w_{s}} + w_{s} \overset{\bullet}{T_{s}} \overset{\circ}{C_{V}} = C_{V} \left(T_{0a} \overset{\bullet}{w_{a}} + T_{0p} \overset{\bullet}{w_{p}} \right)$$
 (26)
$$- \overset{\bullet}{H}_{L} - C_{p} \overset{\bullet}{T_{s}} \overset{\bullet}{w_{n}}$$

where:

$$\dot{C}_V = \frac{d (C_V)}{dP} \times \frac{dP}{dt}$$
 by definition

Solving Equation (26) for the rate of change of system temperature with time (T_s) one may simultaneously solve with Equation (25), which on expansion of terms and appropriate manipulation yields the following equation for mass burning rate:

$$\frac{dw}{dt} = \begin{pmatrix} v_s \hat{p} & Pw_{si} & \hat{H}_L & \gamma T_s \hat{w} & w_s T_s \hat{c}_V & Pw_a & T_{0a} \hat{w}_a \\ \frac{R}{S} & -\frac{R}{S} & \frac{L}{C_V} & +\frac{L}{C_V} & \frac{L}{C_V} & \frac{L}{C_V} & \frac{L}{C_V} \end{pmatrix}$$

$$\begin{bmatrix} P_{11} & PV_{S} & m_{1}R_{u} \\ R_{S} & -(R_{S}w_{S})^{2} \end{bmatrix} \begin{bmatrix} \dot{w}_{T} + \dot{w}_{1} + \dot{w}_{21} - \frac{w_{D1}}{w_{D1}} & PV_{S}w_{S} & \frac{\dot{w}_{T}}{w_{T}} + \frac{\dot{w}_{1}}{w_{2}} + \frac{\dot{w}_{21}}{w_{2}} + \frac$$

APPENDIX E

Equations for Wall Temperature

Computations in WALTEM

Ö

Heat loss is computed by coupling transient conduction in the bomb wall with the convective and radiative transfer to the wall. In the wall, the governing equation is

$$\frac{\delta T_{W}}{\delta t} = \frac{\delta^{2} T_{W}}{\delta \chi^{2}} \cdot c \lambda \tag{23}$$

where T_{W} = temperature field within the wall

a = thermal diffusivity

X = radial distance into the wall

The initial condition is a uniform temperature (set at 298K). The boundary conditions are:

at
$$X = 0$$
: $-K \frac{\delta T}{\delta X} = H_L/A_W$

at
$$X + \infty \frac{\delta T}{\delta X} = 0$$

An explicit centered difference scheme is used for the calculation. At the boundaries the scheme is as follows:

at X = 0:

$$T_0^{n+1} = T_0^n + \frac{\alpha \Delta t}{(\Delta X)^2} \begin{bmatrix} H_L \\ A_W K \end{bmatrix}^2 (\Delta X) - 2T_1 + 2T_2^n$$

where n + 1 is the new time level

n the old time level

T₁ the first interior point

T₂ the second interior point

at back boundary X = N AX:

 $T_{N} = T_{N-1}$

N = number of grid points

APPENDIX F
Capsule Summary of Suboutines

Capsule Summary of Subroutines

ACQUIRE	Subroutine to get file data, allow update and start process
HTLOSS	Subroutine to calculate average heat loss value, average heat transfer coefficient
HTFIT	Subroutine to fit the decay portion of the P-t curve to find heat loss coefficients.
SUMUP	Subroutine to print summary sheet of analysis.
REDUCE	Driver for the differential equation solver.
EVES	Subroutine solves N simultaneous first order differential equations by the Adams method.
SETUP	Subroutine sets the initial conditions for the integration and defines the accuracy required in the solution.
·DIFEQ	This subroutine evaluates the derivative values at each proposed step in the solution. On each call II contains the current time, and Y's are the current integrated values. The results of a call may be rejected and the step size cut in order to maintain accuracy.
PRINT	This subroutine is called to output accepted values during the integration procedure.
FINDTP	A subroutine for table lookup with linear interpolation A direct access read lookup modification of FIND1.
FIND1	A subroutine for table lookup with linear interpolation Extrapolates values out of table range, remembers last argument value and bogins search from that value.
SIMEQ	A subroutine solving L equations in M unknowns, with N sets of right hand constants.
RCALC	Subroutine to do least squares fit on burn rate data.
VPLOT	Versatech* unit plotting package.

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APPENDIX G

Program Listing. CBRED

9823

END

```
MAIN DRIVER FOR LATEST REDUCTION PROGRAM, 10/14/75
1889
             COMMON RUNID(5), RTITLE(18), DATE(3), OPERN(5), PROPN(18),
            1 PSORCE(18), PRLOT(18), PREM(18), TIGHR(18), GAGE(18),
            2 SPACE1(36), CHGWT, WID, PTEMP, WGHTI, PCORR, BYOL, BTEMP,
            3 SCAL, CALCON, STIME, FFID, GL.OD, PD, WOD, F1, DUM, XMUX1, ET1, GAM1.
            4 SPACE2(4). ISK1.KTOT.MZ.MY.PMAX.TMAX.NPMA.PPM.DPMAX.IHL.H.
            5 RHOC.RHO.T10.T90.T1090.P9(5).DP9(5).NPTH.PX(20).
            6 FX(20), ETX(20), XMLX(20), GAMX(20), IDFF, NPA, WEBX(20), ABX(20),
            7 ISK, ISK2, MEM2, XMI, RP, TZERO, VOL, CP, CST, TM, PIGN, TIGN, PTHEO,
            3 CON1.CON9.MEM3.MEM7.SFAZ.L9.P.DP.TSTOP.19.
            9 L8.CONW.TBW.FRAC.Z.SRAT.XTBW.RATE.RL
8882
             COMMON /GENE/ WIDP(5), WODP(5), UCR(5), SFAC(5,5), TDCL(5),
            I WDEY, TDEY, TUST
0003
             COMMON /BLAH/ PLO, PHI, ACO, XNC, RSQ, RMS
8004
             COMMON /KEVIN/ TE(35), XT(35)
0005
             DATA DUMMY WOW 1/
0006
             WRITE (11,197) WOW
0007
       197
            FORMAT (A4)
0000
             REWIND 11
0009
             CALL ACQUIRE
             CALL HTLOSS
0010
0011
             SPACE2(3) = 0.0
0012
             IF (IHL.EQ.2) GO TO 1976
0014
             CALL HTF IT
0015
       1976 CONTINUE
0016
             CALL SUMUP
0017
             CALL REDUCE
0010
             CALL RCALC
0019
             TYPE 100
8020
       188 FORMAT (///)
             CALL VPLOT
0021
0022
             STOP
```

```
1888
            SUBROUTINE ACQUIRE
      C
            SUBROUTINE TO GET FILE DATA, ALLOW UPDATE AND START PROCESS
            COMMON RUNID(5).RTITLE(18).DATE(3).OPERN(5).PROPN(18).
0002
           1 PSORCE(18).PRLOT(18).PREM(18).TIGNR(18).GAGE(18).
           2 SPACE1(36), CHGWT, WID, PTEMP, WGHTI, PCORR, BVOL, BTEMP.
           3 SCAL.CALCON.STIME.FFID.GL.OD.PD.WOD.F1.DUM.XMWX1.ET1.GAM1.
           4 SPACE2(4), ISK1,KTOT,M2,MY,PMAX,TMAX,NPMA,PPM,DPMAX,IHL,H,
           5 RHOC.RHO.T10.T90.T1090.P9(5).DP9(5).NPTH.PX(20).
           6 FX(20), ETX(20), XMUX(20), GAMX(20), IDFF, NPA, WEBX(20), ABX(20),
           7 ISK.ISK2.MEM2.XMI.RP.TZERO.VOL.CP.CST.TM.PIGN.TIGN.PTHEO.
           8 CON1, CON9, MEM3, MEM7, SFAZ, L9, P, DP, TSTOP, 19,
           9 L8, CONW, TBW, FRAC, Z, SRAT, XTBW
            COMMON /GENE/ WIDP(5), WODP(5), UCR(5), SFAC(5,5), TDEL(5),
0003
           1 WDEV, TDEV, TUST
            COMMON /BLAH/ PLO, PHI, ACO, XNC
0004
0005
            DIMENSION A9(4)
            DATA AY/'Y'/.A9/0.25,.375,.50,.625/.BL/
0006
           1. AVG/'AVE '/, P1/'1P '/, P7/'7P '/, P19/'19P '/,
           2 SP/'SPH '/,CY/'CY '/
0007
            DIMENSION LIST (400)
9999
            EQUIVALENCE (LIST(1).RUNID(1))
0009
            EQUIVALENCE (IBUG.LIST(390))
            EQUIVALENCE (ISK1.LIST(399)), (KTOT.LIST(400))
0010
            EQUIVALENCE (ATH, SPACE1(10)), (PBLOW, SPACE1(11)),
0011
            1 (FAID.SPACE1(12)).(T9ID.SPACE1(13)).(CAID.SPACE1(15)).
           2 (XMA.SPACE1(14))
             CALL ASSIGN (2,'SY:CBDAT.001'.0,'SCR')
9912
            DEFINE FILE 2 (3000.6.U.MZ)
0013
9914
            TYPE 300
0015
       300 FORMAT (SX. 'ENTER SMOOTHED TAPE ID'/)
0916
            CALL ASSIGN (3.'FILNAM.EXT',-1.'RDO')
             DEFINE FILE 3(3100.4.U.MY)
9917
9190
            WDEV - 0.0
            TDEV . 0.0
0019
            DO 69 1 - 1.188
0020
             J • 4*1
0051
0022
            L = J - 3
        69 READ (3'1) (LIST(K), K + L.J)
0023
            K . 0
0024
             18UG - 0
0025
             DPMAX . 0.0
0026
6027
             RHOC - SPACE2(1)
            PMAX . 0.0
0020
            FAID - 362000.
0029
            TAID - 3880.
0030
            RU - 1544, *1.8
1500
0032
            X竹 - RU*TAID/FAID
            CAID . 0.0
0033
            ATH - 0.0
0034
            P8LOW - 8.8
0035
0036
            KM - KTOT + 188
0037
              DO 68 J - 101.KM
```

0038	READ (3'J) B1. C1
0039	i = J - 100
0040	II = (I-1)*ISK1 + 1
0041	TI = I1*STIME*1.0E-03
0042	K = K + 1

```
0043
            IF (81.LT.PMAX) GD TO 88
0045
            PMAX = 81
8846
            TMAX - TI
0047
            NPMA - K
0049
            IF (C1.LT.DPMAX) GO TO 68
            PPM - 81
0050
0051
            DPMAX - CI
        68 URITE (2'1) TI, 81, C1
0052
P053
            MEM7 = 1
            PDES - 0.99*PMAX
0054
0055
            CALL FINDTP (PDES.TMAX.3.2.1.2.NPMA.MEM7.MZ)
            PDES = 0.1*PMX
0056
            CALL FINDTP (PDES.T10.3.2.1.2.NPMA.MEM7.MZ)
0057
            PDES = 0.9*PMAX
0050
0050
            CALL FINDTP (PDES. T90.3.2.1.2.NPMA. MEM7. MZ)
             11998 - T98 - T18
0060
1206
            SUM . 0.8
8062
            00 62 I - 1.4
0063
            P9(I) = A9(I)*PMAX
0064
            CALL FINDTP (P9(1).DP9(1).3.2.3.2.NPMA, MEN7.NZ)
0065
        62 SUM * SUM + DPS(I)
            DP9(5) . SUM/4.8
<del>08</del>66
            P9(5) - AVG
0067
0069
            ENDFILE 3
9869
            FRAC - 1.8
6799
            IHL + 2
0871
            H - 0.0
            XMI - 68.62
0972
0073
             IF (PCORR.GT.150000.) XMI = 25.
             TYPE 301
0075
       301 FORMAT (5X.' IS THIS TO BE A'/
0076
            I SX. 'STANDARD ANALYSIS. Y OR N?')
8877
            ACCEPT 302, ANSP
9979
       302 FORMAT (A1)
0079
            CALL PLOT (27.12.4)
0930
            SPACE2(2) - 10.1
6991
            PLO - 0.1
6085
            PHI . 0.8
             IF (ANSP.EQ.AY) GO TO 99
6663
0865
             TYPE 303, RHOC
9899
       303 FORMAT (SX. DENSITY", F12.5)
            ACCEPT 384. TEMP
9697
       384 FORMAT (6E12.8)
6666
             IF (TEMP.NE.8.8) RHOC . TEMP
0089
0091
             TYPE 305. PCORR
0092
       385
            FORMAT (5X. 'IGHITER IMPETUS', F12.5)
6993
            ACCEPT 304. TEMP
            IF (TEMP.NE.B.B) PCORR - TEMP
8894
0896
            TYPE 386. UCHTI
       386
            FORMAT (SX. 'IGNITER LEIGHT' .F12.5)
0097
0096
            ACCEPT 304. TEMP
             IF (TEMP.NE.B.O) WENT! . TEMP
8899
```

0101		IF (PCORR.GT.150000.) XM1 = 25.
0103		TYPE 396, XMI
0184	396	FORMAT (5X, 'IGNITER MOL WEIGHT', F12.5)
0105		ACCEPT 394, TEMP
0106		IF (TEMP.NE.0.0) XMI - TEMP

0166	98	TYPE 323, F1
0167	323	FORMAT (5X, 'PROPELLANT IMPETUS', F12.5)
0168		ACCEPT 304, TEMP
0169		IF (TEMP.NE.0.0) F1 = TEMP
0171		TYPE 324, ET1

15%, 'DIST BURNT - TOTAL BURN AREA')

0231		DO 43 I = 1.NPA
0232	43	ACCEPT 304, WEBX(I), ABX(I)
0233		GO TO 22
0234	21	TYPE 318, GL
0235	318	FORMAT (5%, 'GRAIN LENGTH', F12.5)

ACCEPT 2469, TEMP

TYPE 7354, 18UG

IF (TEMP.NE.0.0) PBLOW - TEMP

7354 FORMAT (5X,'DIFEQ TRACE CONTROL'/

0294

0295

0297

1 5%, TYPE ''1' FOR TRACE ON', [3)
ACCEPT 7355, IBUG
7355 FORMAT ([1])
GO TO 92
99 CONTINUE 0300

and the supplied white products the re-

```
V01-11 SOURCE LISTING
                                                          PAGE 886
  RT-11 FORTRAH IV
              IDFF - 0
  0303
              IF (FFID.EQ.P1) IDFF = 1
  0304
              IF (FFID, EQ, P7) IDFF = 2
  0306
  0308
              IF (FFID.EQ.P19) IDFF = 3
  0310
              IF (FFID.EQ.SP) IDFF - 4
  0312
              IF (FFID.EQ.CY) IDFF = 5
  0314
              IF (IDFF.NE.0) GO TO 87
              TYPE 86
  0316
          86 FORMAT (5X, 'UNKNOWN GRAIN TYPE IN STANDARD'/
  0317
             1 5X, 'PROGRAM HAS ABORTED')
              STOP
  0318
  0319
          87 CONTINUE
              NPTH = 4
  0320
              TEMP1 = XMLX1
  0321
              TEMP2 - GAM1
  0322
  0323
              IF (GAM1.GE.1.0) TEMP2 = 1544.*1.8/778./XMLX1/(GAM1 - 1.)
              IF (XMUX1.LT.1000.) TEMP1 = F1*XMUX1/(1544.*1.8)
  0325
              P2 - 1.0
  0327
              DO 47 I = 1.4
  0328
              PX(I) = P2
  8329
0330 .....
              FX(1) = F1
              ETX(I) = ETI
  0331
              XMUX(1) = TEMP1
  0332
              GAMX(I) - TEMP2
  0333
  0334
              P2 - P2 + 1.8
  0335
          47 CONTINUE
  0336
          92 ISK1 = MAX0(ISK1.1)
              IF (IDFF.EQ.2.OR.IDFF.EQ.3) GO TO 6138
  0337
  0339
              WITH - AMINI (WID. WOD)
  0340
              WOD - UMIN
              WID - WMIN
  0341
  0342
         6138 CONTINUE
              19K - 18
  03.43
              ISK2 * 5
  0344
  0345
              RHO = RHOC*.0361111
  0346
              RETURN
  0347
              END
```

```
0001
            SUBROUTINE HTLOSS
      C
             SUBROUTINE TO CALCULATE HEAT LOSS NUMBER - 2 OPTIONS
0002
            COMMON RUNID (5), RTITLE (18), DATE (3), OPERN (5), PROPN (18),
            1 PSORCE(18), PRLOT(18), PREM(18), TIGNR(18), GAGE(18),
            2 SPACE1(36), CHGWT, WID, PTEMP, WGHTI, PCORR, BYOL, BTEMP,
           3 SCAL, CALCON, STIME, FFID, GL, OD, PD, WOD, F1, DUM, XMWX1, ET1, GAM1,
            4 SPACE2(4), ISK1, KTOT, MZ, MY, PMAX, TMAX, NPMA, PPM, DPMAX, THL, H,
           5 RHOC.RHO.T10.T90.T1090.P9(5).DP9(5).NPTH.PX(20).
           6 FX(20), ETX(20), XMLX(20), GAMX(20), IDFF, NPA, WEBX(20), ABX(20),
             ISK, ISK2, MEM2, XMI, RP, TZERO, VOL, CP, CST, TM, PIGN, TIGN, PTHEO,
           8 CON1.CON9.MEM3.MEM7.SFAC.L9.P.DP.TSTOP.19.
            9 L8.CONW.TBW.FRAC.Z.SRAT.XTBW
            EQUIVALENCE (FAID.SPACE1(12)).(TaID.SPACE1(13)).
0003
            1 (XMA.SPACE1(14)), (CAID.SPACE1(15))
0004
            MEM2 - 1
0005
            CALL FIND! (PMAX.F.PX.FX.NPTH.MEM2)
            CALL FIND: (PMAX.ETA.PX.ETX.NPTH.MEM2)
0006
            CALL FIND: (PMAX,TZP,PX,XMWX,NPTH,MEM2)
0007
0008
            CALL FIND: (PMAX.CVP.PX.GAMX.NPTH.MEM2)
0009
            RU = 1544.*12.*1.8
            XMU = RU*TZP/(12.4F)
0010
            TZERO - TZP
0011
0012
             TA = 298.
            TI + 12. *PCORR*XMI/RU
0013
            VOL - BVOL/(2.54**3)
0014
            CA = 14.7*VOL*29./(298.*RU)
0015
            CI - WGHTI/454.
0016
            CP - CHGUT/454.
0017
0018
            CST - CA + CI
0919
            ETX(20) * CA/29. + CI/XMI + CAID*FRAC/XMA
0020
            TM = CA*298. + CI*TI + CAID*FRAC*TAID
            ETX(19) - CVP*(CST + FRAC*CAID)
0021
0022
            TM - TM/(CST + FRAC*CAID)
            CTOT - CA + CI + CP + CAID
0023
            RA - RU/29.
0024
0025
            RI * RU/XMI
            RP = RU/XML
0026
0927
            RAID - RUXMA
            PTHEG = (CA*RA*298. + CI*RI*TI + CP*RP*TZERO + CAID*RAID*TAID)
0028
            1 /(VOL - CTOT*ETA)/1000.
            TH - THYPMAX/PTHEO
0029
            TL - TI*PMAX/PTHEO
0030
            TLI - TAID*PMAX/PTHEO
0031
0032
            PIGN = (CA*TA/29, + CI*TL/XMI + FRAC*CAID*TL1/XMA)*RU/(VOL ~
            1 (CST + FRAC*CAID)*ETA - CP/RHO - (i. - FRAC)*CAID/RHO)/1000.
            CALL FINDTP (PIGN, TIGN, 3.2, 1, 2, NPMA, MEM2, MZ)
0033
            CON9 - VOL - CP/RHO - (1. - FRAC) *CAID/RHO
0034
0035
             IF (H.EQ.0.0) GO TO 10
```

H = (VOL - CTOT*ETA)*CVF*DPDT/RU*1000.*CTOT/

1 (ETX(20) + CP/MU + (1. - FRAC)*CAID/MA)

IF (FRAC.EQ.1.0) RETURN

10 DPDT . (PTHEO - PMAX)/(TMAX - TIGN)

9837

0039

2841		IF (IHL.NE.2) GO TO 1492
0043		IF (FRAC.EQ.1.0) RETURN
0045		HTL1 = H
0046		GO TO 1493
0047	1492	SPACE1(36) = 450.

```
RT-11 FORTRAN IV
                        V01-11 SOURCE LISTING
                                                         PAGE 902
0048
            TYPE 372, SPACE1(36)
0049
       372 FORMAT (5X, "RYERAGE BOMB WALL TEMPERATURE", F12.5)
0050
            ACCEPT 373, TEMP
       373 FORMAT (E12.0)
0051
0052
            IF (TEMP.NE.8.8) SPACE1(35) . TEMP
0054
            HB - H*1000. SPACEZ(2)/((TZERO*PMAX/PTHED + TM)/2.
           1 - SPACE (36)
0055
            LIB = ((1. - FRAC) *CAID + CP)/(TMAX - TIGN) *1000.
            WB . WB word . B
0056
0057
            H - HB/WB
0050
            TYPE 932, HB. H
0059
       932 FORMAT (2E16.6)
0060
            SPACE1(34) - 8.0
0061
       1493 IF (FRAC.EQ.1.0) GO TO 68
0063
            TST = TIGN - 0.50
            TST - AMAXI(TST, 0.0)
0064
            SUM - 0.0
0065
            DELT - 0.025
0066
0067
            DO 69 I - 1.28
0068
            CALL FINDTP (TST.DP.3.1.3.2.NPMA.MEM2.MZ)
0069
            SUM - SUM + MP
            TST - TST + DELT
0070
        69 CONTINUE
1760
            DPB - SUMKSB.
0072
0973
            PIG - PIGH*1000.
            DYN - FRAC*CAID/TIGN*1000.
0074
            DYNL - DYN
0075
            TSYS - (COM9 - (FRAC+CAID + CST)+ETA)+PIG/ETX(20)/RU
0876
       1498 DYN . DYN . B
9977
9978
            HTL - HTLI
0079
            IF (IHL.EQ.2) GO TO 1494
0091
            HTL . H*DYN*SPACE2(2)*(TSYS - 298.)/1080.
0002
       1494 CONTINUE
            DYN - ((CON9 - (FRAC*CAID + CST)*ETA)*DPB + ETX(28)*RU*HTL/
0003
           1 ETX(19))/(ETX(20)*RU*(TAID - TSYS)/(CST + FRAC*CAID) +
           2 RU*TSYS/XMA - PIG/RHO + PIG*ETA)
0084
            IF (ABS (1. - DYN/DYNL) - 0.001) 310.310.311
0095
       311 DYNL - DYN
            GO TO 1498
9986
0087
       310 CONTINUE
0009 1
            XHIGH = 0.8
0089
            TYPE 1510. WHIGH
       1510 FORMAT (5x. POWER ON IGNITER FLOW', F12.5)
0090
0091
            ACCEPT 1511. TEMP
0092
       1511 FURMAT (E12.8)
            IF (TEMP.NE.0.0) XNIGH - TEMP
0093
            SPACE1(33) - WHIGH
6695
```

0096

0097

6099 0099

0100

PB - PIG***OHIGN

1496 FORMAT (E16.6)

68 COUTTINE

SPACE1(34) - DYN/P8
TYPE 1496. SPACE1(34)

0101 RETURN 0102 END

```
5881
            SUBROUTINE HTF IT
            SUBROUTINE TO FIT DECRY TO FIND HEAT LOSS COEFFICIENT
      C
0002
            COMMON RUNID(5), RTITLE(18), DATE(3), OPERN(5), PROPN(18),
            1 PSORCE(18).PRLOT(18).PREM(18).TIGNR(18).GAGE(18).
           2 SPACE1(36), CHGWT, WID, PTEMP, WGHT1, PCORR, BYOL, BTEMP,
           3 SCAL, CALCON, STIME, FFID, GL, OD, PD, WOD, F1, DUM, XTWX1, ET1, GAM1.
           4 SPACE2(4), ISK1.KTOT, MZ.MY.PMAX.TMAX.NPMA, PPM. DPMAX.IHL.H.
           5 RHOC,RHO,T10,T90,T1090,P9(5),179(5),NPTH,PX(20),
           6 FX(20), ETX(20), XMUX(20), GAMX(20), IDFF, NPA, UEBX(20), ABX(20).
           7 ISK, ISK2.MEM2.XMI.RP, TZERO, VOL, CP, CST, TM, PIGN, TIGN, PTHEO,
           8 CON1, CON9, MEM3, MEM7, SFAZ, L9, P. DP, TSTOP, 19,
           9 L8.CONW.TBW.FRAC.Z.SRAT.XTBW
0003
            COMMON /GENE/ WIDP(5), WODP(5), UCR(5), SFAC(5,5), TDEL(5),
           1 WDEV. TDEV. TUST
0004
            COMMON /BLAH/ PLO, PHI, ACO, XXC
0005
            DIMENSION A1(500.1).81(500.1).D1(1.1)
9996
        28 MEM2 - 1
            MEM3 - 1
0007
            NE - NPMA + 50/ISKI
0008
            INOT - (KTOT - NE) /500 + 1
0009
0010
            CTOT * CST + CP
            RU + 12.*1544.*1.8
6011
            CALL FINDI (PMAX, F. PX, FX, NPTH, NEM2)
0012
            CALL FIND! (PMAX, TZP, PX, XMIX, NPTH, MEM2)
8913
            CALL FINDI (PMAX, CVP, PX, GAMX, NPTH, MEM2)
8814
0015
            MM . RU*TZP/(12.*F)
9916
            80T - (ETX(28) + CP/XTL)*RU
2017
            J * 8
0010
            DO 62 I - NE.KTOT. INOT
0019
            READ (21) TI. P. DP
            CALL FINDI (P. ETA. PX. ETX. NPTH, MEM2)
9020
1500
            DETA - (ETX(HEM2+1) - ETX(HEM2))/(PX(HEM2+1) -
            1 PX(MEM2))
8955
            P . P#1888.
            DP - DP*1000.
6053
0024
            DETA - DETA/1000.
0025
            T - P*(VOL - CTOT*ETA)/BOT
0026
            J - J + 1
0027
            Al(J, i) = Tex4
            BI(J.1) - -DP#(VOL - CYOT*ETA - CTOT*DETA*P)*CVP/RU*1080.
6928
           1 *CTOT/(ETX(20) + CP/X0*U)
        62 CONTINUE
9929
9939
            CALL SIMEO (AL.BI.DI.J.I.I)
            SPACE2(3) = D1(1.1)/SPACE2(2)
0031
4832
            SUM - 8.8
            DO 777 17 - 1.J
6633
0034
        777 SUM - SUM + B1(J.1)**2
6935
            RHS - SORT(SUM/(J - 1))
8835
            UB - (CP/(TMAX - TIGN))*1088.
            ₩8.8 × UB • UB
8837
9939
            TAV - (TZERO**4 + TM**4)/2.
            TG00 - (TZERO + TM)/2. - SPACE1(36)
0039
```

8848	HL = H*LB*TGOO - SPACE2(3)*TAV
0041	SUM1 = 0.0
0042	SUM2 = 0.8
0043	DELT? = (TMAX - TIGN)/1888.
0044	T7 - TIGN

RT-11 FORTR	AN IV VO	1-11 SOURCE	LISTING	PAGE 002
0045 836	CALL FINDTP(T7.	.DPT-3.1.3.2	.3000.MEM2)	
0046	SUM1 - SUM1 + 7	DPT*DELT7		
8847	DPT - AMAXL (DP	T.0.00001)		•
0846	SUM2 - SUM2 + 1	DELT7*DPT***	.8	
9049	77 - 17 + DELT	7	•	
8856	IF (T7.LE.TMAX) GO TO 836		
ยปร2	FAC7 = SUM1**0	.8/SUM2		
0053	H = HL/WB/TG00			
C	H - HWFAC7			
0054	TYPE 100. D1(1.	.1) . SPACE2(3). RMS. H	
0055 100	FORMAT (4E15.5)		••••	
0056	RETURN	•		
0057	END			

```
0001
             SUBROUTINE SUMUP
             SUBROUTINE TO PRINT SUMMARY SHEET OF ANALYSIS
0002
             COMMON RUNID(5).RTITLE(18).DATE(3).OPERN(5).PROPN(18).
            1 PSORCE(18), PRLOT(18), PREM(18), TIGHR(18), GAGE(18),
            2 SPACE1(36), CHGWT, WID, PTEMP, WGHTI, PCORR, BVOL, BTEMP,
            3 SCAL, CALCON, STIME, FFID, GL, OD, PD, WOD, F1, DUM, XMWX1, ET1, GAM1,
              SPACE2(4), ISK1, KTOT, MZ, MY, PMAX, TMAX, HPMA, PPM, DPMAX, IHL, H,
              RHOC, RHO, T10, T90, T1090, P9(5), DP9(5), NPTH, PX(20),
            6 FX(20), ETX(20), XMWX(20), GAMX(20), IDFF, NPA, WEBX(20), ABX(20),
              ISK. ISK2. MEM2. XMI. RP. TZERO. VOL. CP. CST. TM. PIGN. TIGN. PTHEO.
            8 CON1.CON9.MEM3.MEM7.SFAC.L9.P.DP.TSTOP.19.
            9 L8.CONU, TBW, FRAC, Z, SRAT, XTBW
0003
             DIMENSION ABC (5)
             DATA ABC/'CONS', 'TANT', 'PROP', 'ORTI', 'ONAL'/
0004
0005
             PRINT 100
             FORMAT (1H1///)
2006
        100
             PRINT 101. (RUNID(I), I = 1.5)
7000
0008
             FORMAT (10%, 'RUN ID: '29%, 5A4)
9909
             PRINT 102. (RTITLE(I). I = 1.18)
             FORMAT (10X, 'RUN TITLE: ', 25X, 18A4)
0010
0011
             PRINT 103. (DATE(I), I = 1.3)
8012
        103
             FORMAT (10X, 'DATE: ', 30X, 3A4)
9013
             PRINT 104. (OPERN(I). I = 1.5)
            FORMAT (10X, 10PERATOR: 1, 26X, 5A4/)
        184
0014
0015
             PRINT 105
9816
             FORMAT (10X, 'PROPELLANT DATA'')
9917
             PRINT 136, (PROPN(I), I - 1,18)
9919
        186
             FORMAT (10X, 'TYPE: ', 30X, 18A4)
0019
             PRINT 107, CHGWT
0020
        107
             FORMAT (10X, "WEIGHT (GMS): ".22X, F12.5)
             PRINT 100, RHOC
1500
             FORMAT (10X, 'DENSITY (GM/CC): ', 19X, F12.5)
9922
        128
             PRINT 109. PTEMP
0023
0024
        189
             FORMAT (10X.'INITIAL TEMPERATURE (DEG K):'.7X.F12.5)
             PRINT 110. _(PRLOT(I). I = 1.18)
0025
            FORMAT (10X2 LOT: 1,31%, 18A4)
0025
8827
             PRINT [11, (PSORCE(I), [ = 1,18)
9658
        111 FORMAT (10X. 'SOURCE: '. 28X. 18A4)
0029
             PRINT 112, FFID
            FORMAT (10X. 'GRAIN TYPE: '.24X.A4)
0030
0031
             PRINT 113, GL, OD, PD
9932
        113 FORMAT (19X.'LENGTH.OD.ID (IN):'.17X.3(F12.5.'.'))
6633
             PRINT 114. WID. WOD
        114 FORMAT (10x, 'INNER WEB.OUTER WEB (IN):', 10x, 2(F12.5, ', '))
0034
0035
             PRINT 115, FX(1)
1036
            FORMAT (10X, THEORETICAL IMPETUS (FT-LB/LB): 1,4X,F12.5)
0037
             PRINT 116. TZERO
9569
             FORMAT (18X. 'FLAME TEMPERATURE: '. 17X, F12.5)
0639
             PRINT 117, XMLX1
             FORMAT (10X, 'AVERAGE MOLECULAR WEIGHT OF PROD: ', 2X, F12.5)
0040
       117
             PRINT 118, ETX(1)
6041
            FORMAY (18X. 'CO-VOLUME (CU IN/LB): '.14X.F12.5)
0042
```

0043		PRINT 119, GAM1
0044	119	FORMAT (10x. 'GAMMA (RATIO OF SP HTS): '11X.F12.5)
0845		PRINT 120, (PREM(I), I = 1,18)
0046	120	FORMAT (10X, 'REMARKS: ', 27X, 18A4/)
8847		PRINT 121

```
121 FORMAT (10X, 'IGNITER DATA: '/)
0048
0049
            PRINT 122, (TIGNR(I), I = 1.18)
0050
       122 FORMAT (10X, TYPE: 1,30X,18A4)
6651
            PRINT 123, WGHTI
       123 FORMAT (10X, 'WEIGHT (GMS):', 22X, F12.5)
0052
            PRINT 124, PCORR
0053
       124 FORMAT (10X, 'IMPETUS (FT-LB/LB):', 16X, F12.5/)
0054
            PRINT 125
0055
       125 FORMAT (10X, 'EQUIPMENT DATA'/)
0056
            PRINT 126. BVOL
0057
       126 FORMAT (10X, 'BOMB VOLUME (CC):', 18X, F12.5)
0058
            PRINT 127, BTEMP
0059
       127 FORMAT (10X, 'BOMB TEMP (DEG K): '17X, F12.5)
0060
            PRINT 128, (GAGE(I), I = 1,18)
0061
       128 FORMAT (18X, 'GAUGE TYPE: '.24X, 18A4)
0062
            PRINT 129,SCAL
0063
       129 FORMAT (10X, 'CALIBRATION FACTOR (PC/PSI):',7X,F12.5/)
0064
           . PRINT 130
0065
       130 FORMAT (18X, 'RESULTS: '/)
0066
            PRINT 131, PTHEO
0067
       131 FORMAT (10X, 'THEORETICAL MAX PRESS (KPSIA): ',5X,F12.5)
0068
0069
            PRINT 132, PMAX
       132 FORMAT (10X, 'OBSERVED MAX PRESS (KPSIA):'. 8X,F12.5)
0070
            PRINT 133, PIGN
0071
       133 FORMAT (10X, 'IGNITER PRESSURE (KPSIA): '.10X, F12.5)
0072
0073
            PRINT 134
       134 FORMAT (/10X.'IGNITION TIME INFORMATION:')
0074
            PRINT 135, T10
0075
       135 FORMAT (10X. 'TIME TO 10x PMAX (MSEC):'.11X.F12.5)
0076
            PRINT 136. T90
0077
       136 FORMAT (10X. 'TIME TO 90x PMAX (MSEC): '. 11X. F12.5)
9799
            PRINT 137. THAX
0079
       137 FORMAT (10X. TIME TO 100x PMAX (MSEC): 10X.F12.5)
0080
1999
            PRINT 138, T1090
       138 FORMAT (19X. TIME FROM 18% TO 98% PMAX (MSEC): '.2X.F12.5)
0085
0093
            PRINT 148
       148 FORMAT (1H1////18X.
0004
                                          'QUICKNESS INFORMATION: '/)
            PRINT (41, DP9(1)
0005
       141 FORMAT (10X, 'PDOT AT .250 PMAX: '.5X, F12.5)
0086
0087
            PRINT 142, DP9(2)
0088
       142 FORMAT (18X, 'PDOT AT .375 PMAX: ',5X,F12.5)
0089
            PRINT 143. DP9(3)
       143 FORMAT (10X.'PDOT AT .500 PMAX:'.5X.F12.5)
0090
            PRINT 144, DP9(4)
0091
       144 FORMAT (10X.'PDOT AT .625 PMAX: '.5X.F12.5)
0092
            PRINT 145. DP9(5)
6893
       145 FORMAT (10X.'AVERAGE PDOT:'.10X.F12.5/)
0094
            PRINT 139. DPMAX, PPM
0095
       139 FORMAT (10X, MAXIMUM PDOT (MPSI/SEC): Fil.5,5X, OBSERVED AT'.
0096
           1 ' P • (KPSIA):'.F12.5/)
8897
            GO TO (1.2). IHL
           PRINT 146. (ABC(1), 1 * 1.2)
```

0899	146	FORMAT (18%, "HEAT LOSS OPTION: 1,2%,284)
8 1 6 6		GO TO 3
0101	1	PRINT 147, (ABC (I), I = 3.5)
0102	147	FORMAT (10X, 'HEAT LOSS OPTION: ', 2X, 3A4)
8183	3	PRINT 148. H

PAGE 003 V01-11 SOURCE LISTING RI-11 FORTRAN IV

148 FORMAT (10%, 'HEAT LOSS NUMBER: ',F14.7)
RETURN
RND

CALL EVES (JDQ. TPRNT)

ENDFILE 2

ENDFILE 6

RETURN

END

9816

8817 8818

0019

CALL FORM: (0.0. 19.00.PD.WIDP(1).WODP(1).V0.GL.S.UCR(1))

0046

0047		GO TO 39
8048	2	CALL FORM?(0.0.19.OD.PD.WIDP(I).WODP(I).V0.GL.S.UCR(I))
0049		GO TO 39
0050	3	CALL FORM19(0.0.19.0D.PD.WIDP(I).WODP(I).V0.GL.S.UCR(I))
0051		GD TO 39

```
8852
              CALL FORMSP(0.0.19.0D.PD.WIDP(I).WODP(I).V0.GL.S.UCR(I))
0053
0054
              CALL FORMCY(0.0.19.0D.PD.WIDP(I).WODP(I).V8.GL.S.UCR(I))
0055
       39
            SFRC(J,I) = CPRO(I) / (RHO*V0)
            SUM9 - SUM9 + SFAC(J.I)
0056
            SZERO = SZERO + SFAC(J.I)*S
0057
0050
            SIG(J1.3) = AMINI (SIG(J1.3). UCR(I))
0059
        76 CONTINUE
0060
            TDEL(J) = DDW(J)*TDEV
            IF (J.EQ. 1) TDS - TDEL(J)
0061
            TDEL(J) - TDEL(J) - TDS
0063
0064
            CONTINUE
              CONTINUE
0865
        18
            IF (IDU.EQ.1) GO TO 723
0066
            DO 723 I * 1.4
9969
0069
            K = I + I
0070
            DO 723 J - K.S
0071
            IF (UCk(J).GE.UCR(I)) GO TO 723
0073
            A1 + UCR(I)
            A2 - UIDP(I)
0074
0075
            A3 - UODP(I)
0076
            DO 724 L - 1.5
0077
            A4 - SFAC(L. 1)
0078
            SFAC(L.I) = SFAC(L.J)
0079
            SFAC (L.J) - 94
        724 CONTINUE
0080
            UCR(I) + UCR(J)
1869
9892
            WIDP(I) - WIDP(J)
            WODP(I) + WODP(J)
6993
0084
            UCR(J) = AI
0995
            WIDP(J) . AZ
            WODP(J) - A3
0086
0987
       723 CONTINUE
            L9 . 0
9889
            T(1) - TIGH
8889
0090
            IF (IDFF.EQ.0) UCR(I) . WEBX(NPA)
            IF (IDFF.EQ.8) UCR(2) . LEBX(NPA)
0092
            IF (IDFF.EQ.0) UCR(3) - WE8X(NPA)
0894
            IF (IDFF.EQ.0) UCR(4) - LEBX(NPA)
6696
9999
            IF (IDFF.EQ.0) UCR(S) . LEBX(NPA)
0100
            IF (IDFF.EQ.0) SZERÖ * ABX(1)
            NUMB . SUM9
0105
            IF (IDFF.EO.0) NUMB - 1
9193
            PRINT 100. SZERO. NUMB
0105
       188 FORMAT (/10X.'INITIAL SURFACE AREA (SQ IN):'.F12.5/
0105
           1 10x. 'HUMBER OF GRAINS:', 16x.18)
5010
            Y(1) . CST - UGHTI/454.
9110
            Y(2) - UGHTI/454,
9169
            Y(3) . FRAC*CAID
0110
            Y(4) . 8.8
            Y(5) - Y(3)
0111
0112
            Y(6) . 0.8
```

8113	Y(7) + PIGN
	SIG(1,2) - 0.9991
0114	318(1)27 - 0.0001
0115	SIG(2.2) - 0.0001
0116	SIG(3.2) = 0.0001
0117	SIG(4.2) + 8.081

RT-11 FORT	RAN IV V01-11	SOURCE LISTING	PAGE 003
0118	SIG(5.2) = 0.0001		
8119	SIG(5.3) = CAID		
0120	SIG(6,2) - 0.001		
0121	SIG(7.3) - PBLOW		
0122	SPACE1(31) - 0.0		
0123	IF (IHL.EQ.2) RETU	RN	
0 i 25	DN - (LIGHTI/454. +	FRAC*CAID) /TIGN*1080.	
0126	G = H*DN**0.8		
N127	CALL WALTEM (G.T()),298,,TM)	
0128	CALL WALTEM (G. T(1),290,,TM)	
0129	SFAZ = TM		
0130	DQ 1729 I7 = 1,35		
0131 172	9 TE(17) = XT(17)		
0132	RETURN		
0133	END		

V01-11 SOURCE LISTING

PAGE 881

RT-11 FORTRAN IV

WIP - 0.0

0043	DELJ = TIGH
0044	TDLJ - 0.0
0045	DELI = ABX(20)
0045	DO 75 J = 1.JDO
9847	JI = J + 7

60 TO (28,21), IKL

0104	20	DO 1739 I7 = 1.35
0185	1739	XT(17) - TE(17)
0106		G • H*DH**0.8
0197		CALL WALTEM (G.T(2), SFAZ, TSYS)
9108	97	G = H*DH**8.8

```
RT-11 FORTRAN IV
                         V01-11 SOURCE LISTING
                                                          PAGE 003
            ARP = G*(TSYS - XT(1)) / 1000.*SPACE2(2)
0109
0110
            ARP = ARP + SPACE2(3)*(TSYS**4 - XT(1)**4)/1000.*
            1 SPACE2(2)
0111
            GO TO 22
            ARP - H
      21
0112
0113
      22
            CONTINUE
            DY(1) = -Y(1) /USYS*DNH
0114
0115
            DY(2) = -Y(2) \wedge USYS*DNN
0116
            DY(5) = SPACE1(34)*P**SPACE1(33)
0117
            DY(3) = DY(5) - Y(3) \wedge USYS*DNN
            RUT = DY(1) + DY(2) + DY(3) - Y(4) \wedge USYS*DNN
0118
0119
            RMOL = DY(1)/29, + DY(2)/XMI + DY(3)/XMA - Y(4)/USYS*DNN/XMU
            1 -Y(4)*DXMPノ(XMU**2)
0120
            DY(6) * (TSYS*RWT - TAID*DY(5) + WSYS*TSYS*DCVP/CVP
           1 + ARP/CVP + GAM1*TSYS*DNN + P/(12.*R5YS)
           2 *(DY(5)/RHO - WSYS*DETA - ETA*RWT + VSYS*DP/P - VSYS/
           3 WSYS*(RU/RSYS*(RMOL
           4 - TOTMOL*RUT/USYS))))/(TZP - TSYS + P/(12,*RSYS)*
           5 (ETA - 1..7RHO + VSYS/USYS*(1. + RU/RSYS*(1./XMU - TOTMOL/
           6 WSYS))))
0121
            DY1 = DY(5) + DY(6)
            DY(4) = DY(6) - Y(4) / USYS*DNN
0122
0123
            DY(7) = DP
            GO TO (92.93). IHL
8124
0125
        92 IF (ABS(1. - DY1*1000./DN).LE.0.001) GO TO 94
            IF (DYI.LE.O.0) GO TO 93
0127
0129
            DN - DY1*1000.
0130
            GO TO 97
0131
        94 IF (ABS(1. - XT(1)/XT1), LE.0.0001) GO TO 93
0133
            XT1 = XT(1)
0134
            GO TO 20
            IF (Y(8).EQ.0.0) SZERO - AAB
0135
       93
            IF (Y(8).EQ.0.0) ZZERO - Y(6)
0137
            SRAT - AAB/SZERO
0139
            IF (TDEV.NE.0.0) SRAT - SRAT/10.
0140
0142
            2 - (Y(6) - ZZERO)*454./CHGWT
0143
            IF (AAB.GT.0.0) GO TO 675
0145
            RATE . RL
            GO TO 676
0146
0147
       675 RATE - DY(6)/(RHO+AAB)
0148
       676 CONTINUE
0149
            DO 745 I = 1.JDO
0150
            1 - 1 + 7
            DY(J) = RATE*ARATJ(1)
0151
0152
       745 CONTINUE
            RL - RATE
0153
            RATE . RATE*1000.
8154
0155
            FX(20) - TSYS
0156
            RETURN
            END
0157
```

```
SUBROUTINE PRINT (T, Y, DY, N, TPR)
1999
0002
             COMMON RUNID(5), RTITLE(18), DATE(3), OPERN(5), PROPN(18),
            1 PSORCE(18), PRLOT(18), PREM(18), TIGNR(18), GAGE(18),
            2 SPACE1(36), CHGWT, WID, PTEMP, WGHTI, PCORR, BVOL, BTEMP,
            3 SCAL, CALCON, STIME, FFID, GL, OD, PD, WOD, F1, DUM, XMWX1, ET1, GAM1.
            4 SPACE2(4), ISKI, KTOT, MZ, MY, PMAX, TMAX, NPMA, PPM, DPMAX, IHL, H,
            5 RHOC.RHO.T10.T90.T1090,P9(5).DP9(5).NPTH.PX(20).
            6 FX(20), ETX(20), XMUX(20), GAMX(20), IDFF, NPA, WEBX(20), ABX(20),
            7 ISK, ISK2, MEM2, XMI, RP, TZERO, VOL, CP, CST, TM, PIGN, TIGN, PTHEO,
            8 CONI. CON9, MEM3, MEM7, SFAZ, L9, P. DP, TSTOP, I9,
            9 L8, CONW, TBW, FRAC, Z, SRAT, XTBW, RATE, RL
             COMMON /GENE/ WIDP(5), WODP(5), UCR(5), SFAC(5.5), TDEL(5),
0003
            1 WDEV, TDEV, TUST
             COMMON /KEVIN/ TE(35), XT(35)
0004
             DIMENSION T(2), Y(2), DY(2), TPR(2)
0005
             DIMENSION JD(2)
0006
             EQUIVALENCE (JD(1).SPACE1(35))
0007
             DATA XYZ/101000./
8000
0009
             JD0 = JD(1)
             IF (L9.E0.0) L6 - 0
0010
0012
             IF (L0,E0.0) L8 = 0
0014
             IF (L9.EQ.0) L7 = 0
             DO 1729 17 = 1.35
0016
       1729 TE(17) * XT(17)
0017
             IF (TPR(3).E0.5.0) SPACE1(34) = 0.0
0416
0020
             SFAZ - FX(20)
             ISTP - 0
0021
0022
             DO 669 I - 1.JDO
             J = I + 7
0023
             IF (Y(J).LT.UCR(I)) GO TO 669
0024
             ISTP . ISTP + I
0026
0027
        669 CONTINUE
0028
             IF (ISTP.EQ.JDO) N - 0
             IF (L9.E0.0) GO TO 777
0030
             XDO - JDO + 7
9932
             IF (TPR(3).LT.8.0.OR.TPR(3).GT.XDO) GO TO 4936
8033
             ABX(20) * T(1)
0035
       4936 CONTINUE
0036
0037
             \times 00 + 1.0
             IF (TPR(3).EQ.7.9) SPACEI(31) - SPACEI(10)
0038
0040
             IF (TPR(3).NE.XDO) RETURN
0042
        777 CONTINUE
0043
             P - P/1000.
             DP - DP/1000.
0944
0045
             DIST • 2.*Y(8)
             DPRAT - DP/PMAX
0046
0047
             IF (MOD(L9.1SK).NE.0) GO TO 676
             IF (MOD(L7.40).NE.0) GO TO 677
0049
             PRINT 100. (RUNID(1), 1 = 1.5)
0051
0952
       677 L7 . L7 + 1
0053
             IF (MOD(L6.30) NE.0) GO TO 032
0055
             CALL PLOT (27, 12, 4)
```

ť	RT-11	FORT	VI NAS	V01-11	SOURCE I	LISTING	PAGE	002	
	0061		• • • • • • •	L9. ISK2) .NE	.0) GO T	675			
	0063		L8 = L8	+ 1					
	0064		WRITE (1.384) T(1)	, P. DP.	RATE			
N	9965	384	FORMAT (4E13.6)				•	
	0066	675	CONTINUE	,					
	0067		T22 - T0	(1)					
	9068		L9 - L9	+ 1	·			•	
	0069		IF (T)	1).GE.TSTOP) N = 8				
	0071		IF (N.EC	1.0) WRITE (11,384)	xYZ, XYZ, XY	Z,XYZ		
	0073		RETURN						•
	8874	100	FORMAT (1H1/5X,5A4/	7X, TIME	'.6X, 'PRES	S'.6X .'DP/	DT', 8X.	
		``•	ille",	6X21W1 FR12	4X2 SURF	FK - 41. 1	LE BKIL JA.		
			2 'PDOT/F	MAX'/7X, 'MS	EC1.6X,1	KPSIA',4X,	'MPSI/SEC' .:	5X.	
				., 27X, , INCH					
	0075	101	FORMAT (2(4X,F7.3).	4X,F8.3.	4X,F7.3,3(4X,F7,4),4X	,F8.3)	
	0076		END						

END

-		
0055 0056	300	S=2.*E+6.*(OD*(.5*PI3-ALPHA)+PRFD*(BETA+.5*(PI-THETA)))*GRLV=E*GRL
0057		Z=1,-V/V8
0058		RETURN
0059		END

0057		E1-6.*(UU2*RT-1.5*PRFD2*(SIN(THETA)+P13-THETA))
0058		E=E+E1
0059		S=S+2.*E1+36.*PRFD*(PI3-THETA)*GRL
0060		GOTO 300
0061	250	TYPE 251

RT-11 F	ORTR	VI NF	V01-	-11	SOURC	E LIST	ING		PAGE	882
0062 0063 0064 0065 0066		FORMATO STOP V=E*GRL Z=1V/VE RETURN END	GRAIN	GEON	ETRY	YIELD"	NO	OUTER	SLIVERS	.*)

```
RT-11 FORTRAN IV
                        V01-1: SOURCE LISTING
                                                         PAGE 001
            SUBROUTINE FORMSP (DB. J. ODI. PD1. WID. WOD. YO. GL. S. UCR IT)
0001
9992
            DATA PI47.785399/PI/3.141593/
6883
            W - D8*2.8
0004
            DDW = (WPD - WID)/2.
0005
            OD - OD1 + DDW
            OD - AMAX1 (OD, 0.0)
8886
0007
            IF (I.NE.0) GO TO 10
0009
            I = I + 1
0010
            R = UD/2.9
            5 = 4. *P[****2
0011
            V0 - 4.*PI/3.*R**3
0012
            UCRIT - OD
0013
0914
            RETURN
0015
            UCRIT - OD
            IF (W.GT.UCRIT) GO TO 308
0016
8018
            R = (0D - U)/2.
0019
            S = 4.*P[*R**2
0020
           RETURN
0021
      300
            S = 0.0
            RETURN
8022
9023
            FND
```

RT-11	FORTRA	N IV	V01-11	SOURCE	LISTING	PAGE	881
9061 9082 9083 9084 9086 9087 9088 9089 9019 9013 9014		B = SQRT IF (A.NE. ACOS = P' RETURN C = B/A D = ATAN	<pre>/1.570796/, (1 A**2), (0.0) GO TO (2 (C) (C) (0.0) D = D</pre>	10	41593/		

```
0001
             SUBROUTINE WALTEM (F.DTINE.DUM1.DUM2)
0002
             COMMON RUNID(5), RTITLE(18), DATE(3), OPERN(5), PROPN(18),
            1 FSORCE(18),PRLOT(18),PREM(18),TIGHR(18),GAGE(18),
            2 SPACE1(36), CHGWT, WID, PTEMP, WGHTI, PCORR, BVOL, BTEMP,
            3 SCAL, CALCON, STIME, FFID, GL, OD, PD, WOD, F1, DUM, XMWX1, ET1, GAM1,
            4 SPACE2(4), ISK1, KTOT, MZ, MY, PMAX, TMAX, NPMA, PPM, DPMAX, IHL, H,
            5 RHOC.RHO, T10, T90, T1090, P9(5), DP9(5), NPTH, PX(20),
            6 FX(20), ETX(20), XMWX(20), GAMX(20), IDFF, NPA, WEBX(20), ABX(20),
            7 ISK.ISK2.MEM2.XMI.RP.TZERO.VOL.CP.CST.TM.PIGN.TIGN.PTHEO.
            8 CON1, CON9, MEM3, MEM7, SFAZ, L9, P, DP, TSTOP, 19,
            9 L8.CONW.TBW.FRAC.Z.SRAT.XTBW.RATE.RL
0003
             COMMON /KEVIN/ TE(35), XT(35)
             DIMENSION TMA(2), TTA(2), DELX(35), DT(35)
0004
             DATA ISU5/0/, ALFA/.01091/, XK/.0006667/, DELX0/.0009/, FAC/0.3/
0005
0006
             TTA(1) = DUML
             TTA(2) - DUM2
0007
0008
             TMA(1) = 0.0
             TSUM = 0.0
0009
0010
             DTIME = DTINE/1000.
0011
             TMA(2) - DTIME
             IF (ISW5.EQ.0) GO TO 86
0012
             IF (DELTCH - DTIME) 90,90,91
0014
        91 DELTIM - DTIME
0015
             GD TO 92
0016
             DELTIM - DELTCH
0017
        90
0018
             TSUM = TSUM + DELTIM
0019
             IF (TSUM - DTIME) 93,93,95
0020
             TSUM - TSUM - DELTIM
0021
             DELTIM - DTIME - TSUM
             TSUM - TSUM + DELTIM
0022
0023
        93 CONTINUE
0024
             MEM7 - 1
             IF ((TMA(2)-TMA(1)).EQ.0.0) GO TO 1942
0825
             CALL FIND! (TSUM. TG. TMA. TTA. 2. MEM?)
0027
0020
             GO TO 1943
       1942 TG . TTA(1)
0029
0030
             GO TO 1943
0031
       1943 QDOT • F*(TG - XT(1))
0032
             0DOT = 0DOT + SPACE2(3)*(TG**4 - XT(1)**4)
             TZERO - QDOT/XK*2.*DELX0 + XT(2)
0033
9934
             DT(1) * ALFA/(DELX0**2)*(TZERO - 2.*XT(1) + XT(2))
             DO 42 I . 2.31
0035
            CON17 - CONA/(DELX(I) **2)
0036
        42 DT(1) - CON17*(XT(1-1) - CONB*XT(1) + CONC*XT(1+1))
0037
0030
             DO 5 I - 1.31
         5 \times T(1) = X'(1) + DELTIM*DT(1)
8939
0040
            XT(32) * XT(30)
9941
             IF (TSUM - DTIME) 92.74.74
0042
        74 DELTIM - DTIME
0043
            RETURN
0044
        96 ISW6 - I
0045
            DO 21 I . 1.35
```

0046		XT(I) = 298.
0047	21	TE(I) = 298.
0048		TYPE 777, ALFA
0049	777	FORMAT (5X, '80MB THERMAL DIFFUSIVITY', F12.5)
0058		ACCEPT 778, TEMP

END

```
0001
               SUBROUTINE RCALC
         C
               SUBROUTINE TO DO LEAST SQUARES FIT ON BURN RATE DATA
   0002
               COMMON RUNID(5).RTITLE(18).DATE(3).OPERN(5).PFOPN(18).
               1 PSORCE(18).PRLOT(18).PREM(18).TIGNR(18).GAGE(18).
               2 SPACE!(36), CHGWT, WID, PTEMP, WGHTI, PCORR, BYOL, BTEMP.
               3 SCAL.CALCON.STIME.FFID.GL.OD.PD.WOD.F1.DUM.XMWX1.ET1.GAM1.
               4 SPACE2(4), ISK1, KTOT, MZ, MY, PMAX, TMAX, NPMA, PPM, DPMAX, IHL, H,
               5 RHOC.RHO.T10,T90,T1090,P9(5),DP9(5),NPTH,PX(20),
               6 FX(20).ETX(20).XMWX(20).GAMX(20).IDFF.NPA.WEBX(20).A9X(20).
               7 ISK, ISK2, MEM2, XMI, RP, TZERO, VOL, CP, CST, TM, PIGN, TIGN, PTHEO,
               8 CONI.CON9.MEM3.MEM7.SFAC.L9.P1.DP1.TSTOP.I9.
               9 L8.CONW.TBW.FRAC.Z.SRAT.XTBW
   0003
               DIMENSION T(400), P(400), DP(400), R(400)
   0004
               COMMON /BLAH/ PLO. PHI. ACO. XNC.RSQ. RMS
               REWIND 11
   0005
   0006
               DO 669 1 - 1.LB
               READ (11,384) T(1),P(1),DP(1),R(1)
   0007
   0000
               FORMAT (4E13.6)
   0009
          669
               CONTINUE
   0010
               PLL - PLO*PMAX
               PUL - PHI*PMAX
   0811
   0012
               K - 8
   0013
               SUM1 - 0.0
   0014
               SUM2 - 0.0
   0015
               SUM3 - 0.0
   0016
               SUM4 - 0.0
   9017
               SUMS - 0.0
   0018
               DO 735 I - 1.L8
               IF (P(I).LT.PLL) GO TO 735
   0019
   8021
                IF (P(I).GT.PUL) GO TO 735
   6923
               K = K + 1
               DUM1 - ALOG (P(1)*1000.)
   8024
   0025
               DUM2 - ALOG (AMAXI(R(I),0.001))
   0026
               SUM1 - SUM1 + DUM1*DUM2
   0027
               SUM2 - SUM2 + DUMI
   0028
               SUM3 - SUM3 + DUM2
   0029
               SUM4 - SUM4 + DUMI ***2
   0030
               SUMS . SUMS + DUME . STURE
   0031
          735 CONTINUE
   0032
               XONC + (SUM1 - SUM2*SUM3/K)/(SUM4 - (SUM2**2)/K)
   8833
               ACO = EXP (SUM3/K - XNC*SUM2/K)
1. 0034
               RSQ = ((SUM1 - SUM2*SUM3/K)**2)/((SUM4 - (SUM2**2)/K)*
               1 (SUMS -- (SUM3*42)/K))
   0035
               SUM- 0.0
   0036
               K - 0
               DO 733 I - 1.L8
   0037
   0039
                IF (P(1).LT.PLL) GO TO 733
               IF (P(1).GT.PUL) GO TO 733
   0040
   0042
               K \bullet K + 1
               DUM1 - ACO*(P(I)*1080.)**XNC
   0043
               DUM1 = (R(I) - DUMI) \wedge AMAXI(R(I), 0.001)
   0044
               SUM . SUM + DUML**2
   0045
```

0046	CONTINUE	4011M44M		2114188 8
0047	RMS - SQRT	(SUIT (K	_	2114100.0
0048	RETURN			
0049	END			

```
VERSATRAN PLOTTER SUBROUTINE FOR CERED
9991
             SUBROUTINE VPLOT
0002
            COMMON RUNID(5), RTITLE(18), DATE(3), OPERN(5), PROPN(18),
            1 PSORCE(18), PRLOT(18), PREM(18), TIGNR(18), GAGE(18),
           2 SPACE (36), CHGUT, WID, PTEMP, WGHTI, PCORR, BVOL, BTEMP,
           3 SCAL, CALCON, STIME, FFID, GL, OD, PD, WOD, F1, DUM, XMWX1, ET1, GAM1,
            4 SPACE2(4), ISKI, KTOT, MZ, MY, PMAX, TMAX, NPMA, PPM, DPMAX, IHL, H,
           5 RHOC, RHO, T10, T90, T1090, P9(5), DP9(5), NPTH, PX(20),
           6 FX(20), ETX(20), XMUX(20), GAMX(20), IDFF, NPA, UEBX(20), ABX(20),
             ISK, ISK2, MEM2, XMI, RP, TZERO, VOL, CP, CST, TM, PIGN, TIGN, PTHEO,
           8 CON1.CON9.MEM3.MEM7.SFAC.L9.P1.DP1.TSTOP.19.
           9 L8.CONW.TBW.FRAC.Z.SRAT.XTBW
0003
            DIMENSION T(400),P(400),DP(400),R(400)
0004
            COMMON /BLAH/ PLO, PHI, ACO, XNC, RSQ, RMS
0005
            DIMENSION ABC(17)
0006
            DIMENSION BCD (15)
9997
            DIMENSION A(11)
0008
           DIMENSION T9(2), P11(2)
0009
            DIMENSION X1(9), Y1(13)
0010
            DIMENSION TIT(16)
0011
            DATA TIT/'P/PM','AX ','PDOT','
                                                ".'RUN ".'ID: ".'RUN ".
            1 'TITL'.'E: '.'PROP'.' TYP'.'E: '.'GRAI'.'N TY'.'PE: '.
           2 'AT '/, S/9999./
0012
            DATA A/1.0.2.0,4.0.6.0.8.0.10.0.20.0.40.0.60.0.80.0.100.0/
            DATA X1/'TIME'.'-MSE','C '.' '.4HP/PM.'AX '.'1.0',
0013
            1 '10.0'.'100.'/.Y1/"PRES'.'S-KP'.'SIA '.'PDOT'.' MPS'.
                         1,10.1 1,11.0 1,110.01,1RATE1.4H-IN/.1SEC 1/
           2 4HI/SE.'C
            DATA ABC/'THE ','CONS', 'TANT','S IN',' THE', ' EQU', 'ATIO',
0014
            1 'N R '.' = A*'.' P**N'. ARE'. 'A: '.'N: '.'FOR '.'P/PM'.
           2 'AX '.' TO'/
0015
            DATA BCD/'COEF'.'FICI'.'ENT '.'OF D','ETER','MINA',
            1 'TION'.' : '.'PER '.'CENT',' ROO'.'T ME'.'AN E'.'RROR'.' :
0016
            REWIND 11
9917
            DO 683 I - 1.L8
0010
            READ (11.304) T(1).P(1).DP(1).R(1)
0019
       384
            FORMAT (4E13.6)
0020
       683
            CONTINUE
0021
            ENDFILE II
8022
            CALL MODE (8.2.0,1.0,-1.)
0023
            CALL MODE (2.8.0, -0.75.0.875)
0024
            CALL MODE (3.6.0.-0.75.3.9)
            CALL MODE (7.5.0.5.0.5)
0025
            TYPE 1976
0026
       1976 FORMAT (5X. DO YOU WANT TO SUPPRESS THE IGNITION'>
0827
           1 5X. TIME LAG ON THE PRESSURE AND DP/DT PLOT? //
           2 5X. TYPE "S" TO SUPPRESS")
0020
            ACCEPT 1977. STS
0029
       1977 FORMAT (A4)
0038
            IF (STS.NE.STR) GO TO 1978
0032
            DATA STR/'S'/
6633
            19(1) = 10.*[FIX(TIGN/10.)]
0634
            DUMA - 0.0
```

0035	GD TO 1979
0036	1978 T9(1) = 0.0
0037	DUMA = 5.0
0038	1979 T9(2) - TMAX
0039	P11(1) = 0.0

```
0040
            P11(2) - PMAX
0041
            CALL SCAN (T9, P11, -2, 440)
8042
            CALL MODE (-8.DUMI.DUM2.DUM3)
9943
            DUM2 - AMAX1 (DUM2, DUMA)
8844
            CALL MODE (8. DUM1, DUM2, S)
0045
            DUM4 - DUM2
0046
            IMUD - BMUD
0047
            CALL MODE (-9. DUM1.DUM2.DUM3)
0048
            DUM2 = AMAX1(DUM2,5.8)
0049
            CALL MODE (9.0.0.DUM2.5)
0050
            CALL AXES (9.2,X1(1),11.2,Y1(1))
0051
            CALL DRAW (T. P. LB. 440)
            CALL SCAN (T. DP. -LB. 440)
0052
0953
            CALL MODE (8.DUMB.DUM4.S)
0054
            CALL MODE (-2.DUM1.DUM2.DUM3)
0055
            DUMS - DUM3 - 0.5
0056
            CALL MODE (2, S.S.DUM5)
            CALL MODE (9, 0.0, S, S)
0057
0058
            CALL AXES (-4.0,X1(4),13.3,Y1(4))
0059
            CALL MODE (2.S.S.DUM3)
            CALL DRAW (T. DP. LB. 448)
0060
            K - 0
0061
            CALL MODE (4,.073,.055,S)
0062
            CALL MODE (6.5..080.5)
8063
0064
            CALL MODE (3,5,5,-.2)
0065
            CALL HOTE (0.0,3,0,TIT(5),7)
9966
            CALL NOTE (1.0.3.0.RUNID(1).20)
            CALL NOTE (0.0.2.8.TIT(7).10)
9067
0060
            CALL NOTE (1.0.2.8.RTITLE(1).68)
0069
            CALL NOTE (0.0.2.6.TIT(10).10)
            CALL NOTE (1.0.2.6.PROPN(1).60)
0970
0071
            CALL HOTE (0.0.2.4.TIT(13).11)
            CALL NOTE (1.0.2.4.FFID.4) -
0072
6073
            IF (K.EQ.0) GO TO 10
0075
            CALL NOTE (0.0.2.2.TIT(3).4)
0876
            CALL NOTE (2.0.2.2.TIT(16).2)
            CALL NOTE (4.0.2.2.TIT(1).6)
9977
0070
            Y + 2.2 - 0.2
0079
            DO 61 1 * 1.4
            CALL HOTE (0.0.Y.DP9(1).1003)
9989
            CALL NOTE (4.0.Y.P9(1).1003)
1866
           Y . Y - 0.2
6685
        61
        10 CALL MODE (4..1..067.5)
0003
6664
            CALL MODE (6.5..1.5)
0085
            CALL MODE (3.S.S.3.8)
9966
            CALL DRAU (0..0..1.9880)
            PMRX - 0.8
8887
            DO 65 1 . 1.L8
6666
9889
        65 PMAX - AMAXI (PMAX. P(1);
6699
            DO 69 I -1.L8
1689
        69 T(1) • P(1) /PMAX
            CALL MODE (8. 8.0, 8.2, 3.0)
8892
```

0093	CALL AXES (6.2,X1(5),13.3,Y1(4))
0094	CALL DRAW (T. DP. L8, 440)
0095	CALL FORM (5, 1.0, 5, 1.8)
0096	K = 1
0097	CALL MODE (4073055.5)

```
0098
            CALL MODE (6,5,080,5)
0099
            CALL MODE (3.5.5.-.2)
0100
            CALL NOTE (0.0.3.0.TIT(5),7)
            CALL NOTE (1.0.3.0.RUNID(1).20)
0101
            CALL NOTE (0.0.2.8.TIT(7).10)
0102
0103
            CALL NOTE (1.0.2.8.RTITLE(1).68)
            CALL NOTE (0.0.2.6.TIT(10),10)
0104
0105
            CALL NOTE (1.8,2.6,PROPN(1),68)
6186
            CALL NOTE (0.0.2.4, TIT(13), 11)
0107
            CALL NOTE (1.0.2.4,FFID.4)
0108
            IF (K.EQ.8) GO TO 11
            CALL NOTE (0.0.2.2.TIT(3).4)
0110
            CALL NOTE (2.0.2.2.TIT(16).2)
0111
0112
            CALL NOTE (4.0.2.2.TIT(1).6)
            Y = 2.2 - 0.2
0113
            DO 91 I - 1.4
0114
            CALL NOTE (0.0, Y, DP9(1), 1003)
0115
            CALL NOTE (4.0, Y, P9(1), 1003)
0116
        91 Y = Y - 0.2
0117
0110
            CALL MODE (4,.1,.067,S)
            CALL MODE (6.5..1.5)
0119
0129
            CALL MODE (3.5.5.3.8)
0121
            CALL DRAW (0..0..1.9000)
            PFAC - 1.0
0155
            IF (PMAX.LE.10.0) PFAC - 10.0
0123
0125
            PMIN = 1.0
            PMIN - PMIN/PFAC
0126
            1G0 - 0
0127
9129
            PSTO . PMAX
            TYPE 529
0129
            FORMAT (5X.'DO YOU WANT ALL THE RATE CURVE?'/
       529
0130
           1 5X. 'ENTER 1 FOR YES. PLEASE')
0131
            ACCEPT 530, IGO
       538 FORMAT (11)
0132
            IF (IGO.EQ.1) GO TO 533
6133
            PMIN . PLO*PMAX
0135
            PSTO - PHI*PMAX
0136
       533
0137
            CONTINUE
8118
            RMAX - 10.0
0139
            RMIN - 0.1
6148
            J97 - 8
            DO 68 1 - 1.L8
0141
            IF (P(I).LT.PMIN) GO TO 68
8142
8144
            IF (P(I).GT.PSTO) GO TO 68
0146
            197 - 197 + L
            P(J97) - ALOG18 (AMAX1(P(I), PMIN))
8147
            R(J97) = ALOGIO (AMAXI(AMINI(R(I).10.0), 0.1))
0148
0149
            CONTINUE
0150
            DUM! - 0.0
            IF (PFAC.GT.1.0) DUML - -1.8
0151
            CALL MODE (0.DUMI. 0.4.
                                       S)
0153
            CALL MODE (9. -1.0. 0.4.
6154
```

0155		DO 66 I = 1.10
0156	66	A(I) = 2.5*ALOG18(A(I+1)/A(I))
0157		CALL FORM (1010.A.1010.A)
0158		IF (PFAC.EQ.1.0) GO TO 997
0160		CALL NOTE (-0.1,-0.25,Y1(8),3)

PAGE 004

0214		CALL NOTE (0.0,Y,DP9(1),1003)
0215		CALL NOTE (4.0, Y.P9(1).1003)
0216	91	Y = Y - 0.2
0217	12	CALL MODE (4,.1,.067.8)
0218		CALL MODE (F.S1.S)

RY-11	FOR TRAN IV	V01-11 S0	URCE LISTING	PAGE 005
0219 0220 0221 0222 0223	CALL DRAW	(3,5,5,3,8) (0,0,1,900 (0,0,0,0,0		

SUBROUTINE FINDTF (ARG. ANS. NVAR, NARG, NANS. IDEV. NPTS.I.MG)
SUBROUTINE FIND. TABLE LOOK-UP WITH LINEAR INTERPOLATION.
FINDTP IS A DIRECT ACCESS READ LOOKUP MODIFICATION OF
FIND. INSTEAD OF AN ARRAY, PASSING THE ARGUMENTS. IT
EXPECTS A FILE WHICH HAS BEEN FILLED WITH NVAR VARIABLES.
TO THE EXTENT OF NPTS RECORDS. THE INDEPENDENT
VARIABLE MAY OCCUPY ANY OF THE LOCATIONS 1-NVAR AND
NARG GIVES ITS LOCATION. LIKEWISE FOR THE DEPENDENT
VARIABLE WHOSE LOCATION IS GIVEN BY NANS. IDEV GIVES
THE DEVICE NUMBER OR FILE NUMBER WHICH WAS DEFINED
IN SETTING UP THE FILE. ALL OTHER FIND COMMENTS
APPLY TO FINDTP.
NOTE.... LIKE FIND. THE INDEPENDENT VARIABLE MUST
BE MONOTONICALLY INCREASING....

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J. ANDERSON 6-1-65

EXTRAPOLATES FOR VALUES OUT OF TABLE RANGE REMEMBERS LAST ARG VALUE AND BEGINS SEARCH FROM THAT VALUE CALLING SEQUENCE IS

CALL FINDTP (ARG.ANS.NVAR.NARG.NANS.IDEV.NPTS.I)

ARG IS THE ARGUMENT
ANS CONTAINS RESULT ON EXIT
X IS ONE DIMENSIONAL ARRAY OF INDEP. VARIABLE
X IS ONE DIMENSIONAL ARRAY OF DEP. VARIABLE
NPTS IS NUMBER OF TABLE ENTRIES
MEM SHOULD BE INITIALIZED TO 1 - AFTER FIRST
CALL THE SUBROUTINE WILL PRESERVE VALUES IN IT

VARIABLE WHOSE LOCATION IS GIVEN BY NAME. IDEV GIVES THE DEVICE NUMBER OR FILE NUMBER WHICH WAS DEFINED IN SETTING UP THE FILE. ALL OTHER FIND COMMENTS APPLY TO FINDTP.

NOTE.... LIKE FIND, THE INDEPENDENT VARIABLE MUST BE MONOTONICALLY INCREASING....

J. ANDERSON 6-1-65

EXTRAPOLATES FOR VALUES OUT OF TABLE RANGE REMEMBERS LAST AND VALUE AND BEGINS SEARCH FROM THAT VALUE CALLING SEQUENCE 'S

CALL FINDTP (ARG.ANS.NVAR.NARG.NANS.IDEV.NPTS.I)

ARG IS THE ARGUMENT
ANS CONTAINS RESULT ON EXIT
X IS ONE DIMENSIONAL ARRAY OF INDEP. VARIABLE
X IS ONE DIMENSIONAL ARRAY OF DEP. VARIABLE
NPTS IS NUMBER OF TABLE ENTRIES
MEM SHOULD BE INITIALIZED TO 1 - AFTER FIRST
CALL THE SUBROUTINE WILL PRESERVE VALUES IN IT

DIMENSION X(10), Y(10)

```
0003 1 READ (IDEV'I) (X(J), J = 1,NVAR)
0004 IF (X(NARG)-ARG) 4,2,2
0005 2 I=I-1
0006 IF (I-1) 3,3,1
0007 3 I=1
```

```
RT-11 FOR TRAN IV
                        V01-11 SOURCE LISTING
                                                         PAGE 002
0006
            GO TO 7
0009
         4 K = 1 + 1
            READ (IDEV'K) (Y(J), J = 1.NVAR)
0010
0011
            IF (Y(NARG)-ARG) 5.7.7
0012
          5 I=I+1
0013
            IF (I-NPTS)
                           4.6.6
0014
          6 I=NPTS-1
         7 K = I + 1
0016
            READ (IDEV'I) (X(J), J = 1, NVAR)
            READ (IDEV'K) (Y(J), J = 1, NVAR)
0017
            ANS=X(NANS)+(Y(NANS)-X(NANS))*(ARG-X(NARG))/(Y(NARG)-X(NARG))
0016
0019
            RETURN
0020
            END
```

```
0001
            SUBROUTINE FIND! (ARG. ANS. X.Y. NPTS. I)
            SUBROUTINE FIND. TABLE LOOK-UP WITH LINEAR INTERPOLATION.
      C
             J. ANDERSON
                              6-1-65
      C
                    EXTRAPOLATES FOR VALUES OUT OF TABLE RANGE
      C
                    REMEMBERS LAST ARG VALUE AND BEGINS SEARCH FROM THAT VALUE
                    CALLIN SEQUENCE IS .....
      C
                       CALL FIND (ARG.ANS.X.Y.NPTS.MEM)
                           ARG IS THE ARGUMENT
                           ANS CONTAINS RESULT ON EXIT
                           X IS ONE DIMENSIONAL ARRAY OF INDEF. VARIABLE
                           X IS ONE DIMENSIONAL ARRAY OF DEP. VARIABLE
                           NPTS IS NUMBER OF TABLE ENTRIES
      C
                           MEM SHOULD BE INITIALIZED TO 1 - AFTER FIRST
      C
                             CALL THE SUBROUTINE WILL PRESERVE VALUES IN IT
            DIMENSION X(10),Y(10)
0002
0003
          1 IF (X(I)-ARG)
                             4,2,2
0004
          2 I=I-1
0005
            IF (I-1) 3.3.1
0006
          3 I-1
            GO TO 7
0007
          4 IF (X(I+1)-ARG) 5,7,7
0008
6009
          5 I=I+1
0010
            IF (I-NPTS)
                           4.6.6
0011
          6 I-NPTS-1
          7 ANS=Y(I)+(Y(I+1)-Y(I))*(ARG-X(I)).*(X(i+1)-X(I))
0012
0013
            RETURN
            END
0014
```

```
0001
            SUBROUTINE SIMEQ (A.B.D.L.M.N)
            SUBROUTINE SOLVING L EQUATIONS IN M UNKNOWNS WITH N SETS OF RIGHT-
            HAND CONSTANTS. A(L.M) IS THE MATRIX OF COEFFICIENTS AND B(L.N) IS
            THE MATRIX OF COLUMNS OF ANSWERS, ON RETURN FROM SUBROUTINE, A
            CONTAINS ORTHOGONALIZED COLUMNS. B CONTAINS THE RESIDUALS. AND
            D(M, N) CONTAINS THE SOLUTIONS, FOR MORE EQUATIONS THAN UNKNOWNS
      C
            THE LEAST-SQUARES SOLUTION IS OBTAINED, OTHERWISE THE SOLUTION IS
      C
            IN TERMS OF THE FIRST L LINEARLY INDEPENDENT VARIABLES. REQUIRES..
            DIMENSION A(15,15), B(15,15), D(15,15)
            CALL SIMEQ (A. B. D. L. M. N)
0002
            DIMENSION A(500.1), B(500.1), C(3.3), D(1.1)
0003
            DO 701 I=1,M
            DO 700 J-1.M
0004
0005
       700
            C(I,J) = 0.0
            DO 701 K=1.N
0006
       701
0007
            D(I,K) = 0.0
            DO 702 J-2.M
9999
0009
            DO 702 I-1.L
0010
       702 \ C(J,J-1) = C(J,J-1) + (A(I,J)*A(I,J))
0011
            DO 712 K-1.M
            DO 709 J-K.M
0012
0013
            DO 703 II=1,L
0014
            C(K,J) = C(K,J) + (A(II,K)*A(II,J))
            IF (K-J) 706, 704, 704
0015
0016
       704
            IF (K-1) 709, 709, 705
            IF (1.E-7*C(K,K-1) - 1.E7*C(K,K)) 789, 712, 712
0017
       705
9190
       706
            C(K,J) = C(K,J)/C(K,K)
            DO 700 12-1.L
0019
       787
            A(12,J) = A(12,J) - (A(12,K)*C(K,J))
0020
       708
0021
       709
            CONTINUE
0022
            DO 711 J2=1.N
0023
            DO 710 13-1.L
0024
       710 D(K.J2) * D(K.J2) + (A(I3.K)*B(I3.J2)/C(K.K))
0025
            DO 711 14-1.L
0026
       711 B(14,J2) = B(14,J2) - (A(14,K)*D(K,J2))
       712 CONTINUE
0027
            IF (M - 1) 715, 715, 714
0020
0029
       714 DO 713 I=2.M
            IT = M+1-1
0030
            JT - IT+1
0031
            DO 713 J+1.N
9032
0033
            DO 713 K-JT.M
0034
       713 D(IT_*J) = D(IT_*J) - (C(IT_*K)*D(K_*J))
0035
       715 RETURN
0036
            END
```

FIGURES IF THE ABSOLUTE VALUE OF THE INTEGRAL IS LESS THAN THIS VALUE. ITS USE IS TO PREVENT USE OF AN INORDINATELY SMALL STEP SIZE AS AN

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INSETUP (T.Y.SIG.N)
DIMENSION T(2).Y(N).SIG(N.3)

THIS SUBROUTINE SETS THE INITIAL CONDITIONS FOR THE INTEGRATION AND DEFINES THE ACCURACY REQUIRED IN THE SOLUTION.

T(1) - STARTING TIME (ASSUMED 8.0 IF NOT SPECIFIED)

T(2) - INITIAL TIME INCREMENT (ASSUMED 1.0E-5 IF UNSPECIFIED)

Y(I) - INITIAL VALUES OF DEPENDENT VARIABLES, I-I.N (ASSUMED 0.0 IF UNSPECIFIED)

SIG(1.1) - REQUIRED ACCURACY FOR THE DEPENDENT VARIABLES.
I+1.N WHERE SIG+1.E-M INDICATES M SIGNIFICANT
FIGURES. (ASSUMED 1.E-3 IF UNSPECIFIED)

C	SIG(1.2) =	MINIMUM ABSOLUTE ACCURACY DESIRED-SIG(1.2)*SIG(1.1
C	•	>. THIS SUSPENDS THE REQUIRED NO. OF SIGNIFICANT
C		FIGURES IF THE ABSOLUTE VALUE OF THE INTEGRAL IS
C		LESS THAN THIS VALUE. ITS USE IS TO PREVENT USE
C		OF AN INORDINATELY SMALL STEP SIZE AS AN INTEGRAL

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V01-11 SOURCE LISTING **PAGE 802** LEAVES 0. (ASSUMED 0. IF UNSPECIFIED) SIG(1.3) - THRESHOLD VALUE FOR THE VARIABLE I. EVES WILL HIT THIS VALUE EXACTLY DURING THE INTEGRATION AND LET THE USER ROUTINES KNOW THIS VALUE HAS BEEN HIT. (ASSUMED 1.0E+35 IF UNSPECIFIED) SUBROUTINE DIFEC (T, Y, DY, N, TPR) DIMENSION T(2), Y(N), DY(N), TPR(2) THIS SUBROUTINE EVALUATES THE DERIVATIVE VALUES (DY(I)).AT EACH PROPOSED STEP IN THE SOLUTION, ON EACH CALL, T(1) CONTAINS THE CURRENT TIME, AND THE Y(1)'S ARE THE CURRENT INTEGRATED VALUES. THE RESULTS OF A CALL MAY BE REJECTED, AND THE STEP SIZE CUT IN ORDER TO MAINTAIN ACCURACY, SO NO PERMANENT SWITCH SETTING SHOULD BE MADE IN THIS ROUTINE, BASED ON PROPOSED VALUES -SUBROUTINE PRINT (T.Y.DY.N.TPR) DIMENSION T(2), Y(N), DY(N), TPR(N+4) THIS SUBROUTINE IS CALLED TO OUTPUT ACCEPTED VALUES DURING THE INTEGRATION PROCEEDURE. THE FREQUENCY WITH WHICH IT IS CALLED FOR VARIABLES SET IN SUBROUTINE SETUP AS FOLLOWS --THE TPRNT ARRAY CONSISTS OF PAIRS OF VALUES DELTA T AND TLIM

IS DEPENDENT ON THE NUMBER OF INTEGRATION STEPS. THE TPRNT ARRAY SUPPLIED TO EVES BY THE MAIN PROGRAM, AND THE THRESHOLD VALUES

SUCH THAT DELTA T IS THE PRINT INTERVAL UNTIL TLIM IS REACHED . WHEREUPON THE NEXT PAIR OF VALUES TAKES CONTROL.

IF DELTA T .EQ. Ø EYERY ACCEPTED POINT IS PRINTED IF DELTA T .GT. 8 PRINTING OCCURS ONLY AT INTERVALS OF DELTA T

IF DELTA T .LT. 9 PRINTING OCCURS AT EVERY ACCEPTED POINT, BUT AMONG THESE POINTS ARE INTERVALS OF ABS(DELTA T)

-REGARDLESS OF THE TPRNT CONTROLS.A CALL TO PRINT WILL OCCUR EACH TIME A VARIABLE REACHES IT'S THRESHOLD VALUE.

THE TPR ARRAY HAS THE FOLLOWING INFORMATION WHEN PRINT IS CALLED

TPR(1) CURRENT DELTA T FROM TPRNT ARRAY

CURRENT TI IM FROM TPRHT ARRAY

8 IF CALLED THRU ACCURACY CONTROLLED STEP SIZE. +J IF CALLED DUE TO VARIABLE J RISING TO ITS

THRESHOLD VALUE

-J IF CALLED DUE TO VARIABLE J FALLING TO ITS THRESHOLD VALUE

N+1 IF CALLED AT SPECIFIED PRINT POINT TPR(4) CURRENT TIME STEP SIZE FOR INTEGRATION (IF THIS VALUE IS CHANGED IN PRINT, THE DIFFERENCE TABLE WILL BE ZEROED AND THE SOLUTION RESTARTED WITH THE ALTERED VALUE AS THE STEP SIZE. THIS IS TO GET AROUND EXTERNALLY

С	INDUCED DISCONTINUITIES WITHOUT REQUIRING THE INTEGRATION
С	TO DISCOVER THE POINT OF OCCURRENCE BY HALVING ITS STEP.)
C	TPR(5) ROUGHEST VARIABLE OF CURRENT STEP
С	TPR(6)-TPR(N+5) CURRENT THRESHOLD SETTINGS FOR EACH VARIABLE
C	(ALLOWING DYNAMIC RESET OF THRESHOLDS)

```
C
      C
0002
            DIMENSION D1(12,7),T1(2),D2(12,7),T2(2),Y1(12),DY1(12),SIG(12,3),
           1 Y2(12), DY2(12), TPRNT(6), TPR(17)
0003
            DIMENSION T3(2)
6004
            EQUIVALENCE (D1(1),Y1(1)),(D1(13),DY1(1)),(D2(1),Y2(1))
            1 (D2(13),DY2(1))
0005
            TBIAS-0.0
            KTERR-0
0006
9007
          1 MaN
          2 DO
0000
                  4 I=1.M
                  3 J-1.7
0009
            DO
            D1(I,J)-0.0
0010
1100
          3 D2(I,J)=0.0
0012
            SIG(I,1)=1.0E-03
9913
            SIG(1.3) =1.0E+35
9914
          4 SIG(1,2) +0.0
0015
             IPRNT-1
0016
            T1(1)=0.0
0017
            T1(2)=1.E-5
      C
             INITIAL SETUP COMPLETE. CALL IN THE PROGRAMMERS SETUP.
      C
            CALL SETUP (TI, YI, SIG, M)
0018
0019
            IF (M.LT.0) IPRNT - -M
            1: " N
0021
9922
            TPR(1) *TPRNT(IPRNT)
0023
            TPR(2)=TPRNT(IPRNT+1)
            THEXT -T1(1)+ABS(TPR(1))
0034
0025
            ISUV-I
0026
            TPR(3) =0.
            DO 5 1-1.M
0027
            J*1+5
0028
          5 TPR(J) *SIG(1.3)
0029
0030
          6 CALL DIFEC(TI.YI.DYI.M.TPR)
0631
             IF (M) 64, 64, 7
          7 CALL PRINT(TL.YL.DYL.M.TPR)
0032
             IF (M) 64, 64, 0
0033
          8 T2(1)*T1(1)*T1(2)
9934
9935
            T2(2) -T1(2)
      C
            PREDICT AHEAD BY ADAMS METHOD.
      C
9936
                  9 I-1.M
            Y2(1) =0.34861111*D1(1.6) +0.375*D1(1.5) +0.41666667*D1(1.4)+0.5*
0037
            1 D1(1.3)+T1(2)*DY1(1)+Y1(1)
0030
             T3(1) *T2(1) +TB1AS
0239
            73(2)-T2(2)
      C
      ¢
            OBTAIN THE DERIVATIVES AT THE PROPOSED POINT.
      C
            CALL DIFEO(T3.Y2.DY2.M.TPR)
0040
```

```
RT-11 FORTRAN IV
                       V01-11 SOURCE LISTING
                                                          PAGE 004
0044
            DO 18 I=1.M
0045
            D2(I,3) =T2(2) *(DY2(I)-DY1(I))
0046
            DO 11 J=3.6
0047
         11 D2(I,J+1) =D2(I,J) -D1(I,J)
      C
      C
            DELETE DIVERGENT DIFFERENCES.
      C
0048
            DO 15 J=3.5
0049
            IF (ABS(D2(I,J+1))-ABS(D2(I,J))) 15, 12, 12
         12 IF (ABS(D2(I,J+2))-ABS(D2(I,J+1))) 15, 13, 13
0050
0051
         13 DO 14 KaJ.6
0052
         14 D2([,K+1)=0.
0053
            GO TO 16
         15 CONTINUE
0054
0055
            J=7
0056
         16 XYZ=0.3*D2(I,J)/(SIG(I,1)*AMAX1(ABS(Y2(I)),1.8E-30,SIG(I,2)))
0057
            IF (ABS(XYZ)-EMAX) 18, 17, 17
0058
         17 EMAX=ABS(XYZ)
0059
            JCRAP . I
0060
            TPR (5) = 1
0061
         18 CONTINUE
      C
      C
            DETERMINE IF ERROR IS WITHIN BOUNDS.
      C
0062
            IF (CMAX-1.0) 30, 30, 19
      C
            ERROR TOO BIG. HALVE THE INTERVAL AND TRY AGAIN.
      C
      C
0063
         19 T1(2)=.5*T1(2)
      C
      C
            CHECK FOR DELTA TIME INSIGNIFICANT.
      C
            IF (T1(1)/T1(2)-1.0E06) 27, 20, 20
0064
0065
         20 CONTINUE
      C
            DELTA TIME IS INSIGNIFICANT. TRANSLATE THE ORIGIN.
            DO 21 IJAZZ=1,M
0066
         21 Y1(IJAZZ) = Y2(IJAZZ)
0067
0068
            TBIAS=TBIAS+T1(1)+2.0%T1(2)
0069
            T2(2) =0.
            T1(1)=0.0
0070
0071
            KTERR=KTERR+1
            IF (KTERR-4) 23, 22, 22
0072
        22 TYPE 67, JCRAP
0073
            RETURN
0074
3075
        23 TYPE 68, KTERR, JCRAP
      C
            RESTART THE SOLUTION.
0076
         24 DO 26 LJAZZ=1.M
```

0077	DO 25 JAZZ=2.6
0078	25 D1(IJAZZ.JAZZ) =0.
0079	DO 26 JAZZ*2,7
0080	26 D2(IJAZZ.JAZZ)=0.
0081	T3(1)=TBIAS+T1(1)

ERROR TOO SMALL. ACCEPT THIS POINT. BUT DOUBLE THE
INTERVAL FOR THE NEXT POINT.

2134

39 T1(2)=2.0*T2(2)

```
0135
            G0 T0 41
      C
            ERROR O.K. BUY THIS POINT AND MAINTAIN CURRENT INTERVAL.
0136
         40 T1(2) =T2(2)
         41 IF (TPR(3) .NE. 0.0) GO TO 73
0137
            IF (TPR(1)) 42, 44, 45
0139
0140
         42 T3(1) =T2(1) +TBIAS
            T3(2) =T2(2)
0141
            OUTPUT THE ACCEPTED POINT.
0142
            TPR(4)=T3(2)
0143
            IF (ABS((T3(1)-TNEXT)/TNEXT) .LE. 0.00001) TPR(3)=M+1
0145
            TZSAVE=TPR(3)
0146
            CALL PRINT(T3, Y2, DY2, M, TPR)
0147
            ISWV=1
0148
            IF (M) 64, 64, 43
      C
            TEST FOR PRINTING CONDITIONS.
0149
        43 IF (T3(2) ,NE, TPR(4)) GO TO 200
0151
            IF (ABS((T3(1)-TNEXT)/TNEXT)-0.00001) 47, 51, 51
0152
         44 TNEXT=TPR(2)
         73 IF (T3(1) - TPR(2)) 46, 45, 46
0153
         45 IF (AB$((T3(1)-THEXT)/THEXT)-0.00001) 99, 51, 51
0154
        99 TPR(3) = M + 1
0155
            TZSAVE = TPR (3)
0156
        46 T3(2) = T2(2)
0157
            T3(2) *T2(2)
0158
3150
            TPR(4)=T3(2)
0160
            CALL PRINT(T3, Y2, DY2, M, TPR)
            ISUV-1
0161
            IF (M) 64, 64, 47
0162
        47 IF (T3(2) .NE. TPR(4)) GO TO 200
0163
            IF (TZSAVE .EQ. TPR(3)) TNEXT=TNEXT+ABS(TPR(1))
0165
            [F (T3(1)~TPR(2)+0.1*TPR(1)) 51, 48, 48
0167
         48 IPRNT=IPRNT+2
0168
            TPR(1) = TPRNT(IPRNT)
0169
0170
            TPR(2) "TPRNT(IPRHT+1)
            IF (TPR(1)) 49, 50, 49
0171
         49 TNEXT = T3(1) + ABS(TPR(1))
0172
0173
            GO TO 51
8174
         50 THEXT TPR(2)
0175
         51 IF (T1(2)+T3(1)-TNEXT) 53, 53, 52
0176
         52 T1(2) "TNEXT-T3(1)
0177
         53 CONTINUE
0178
         54 T1(1)=T2(1)
0179
         55 W-T1(2)/T2(2)
0190
            IF (W-1.0) 56, 65, 56
            ADJUST THE DIFFERENCE TABLE FOR INTERVAL CHANGE.
```

	U					
0181		56	DO	57	I=1.M	
0182				=W*l		
0183			TMP	3 + D:	2(1,3)	
0184			TMP	4-D	2(1.4)	

and alone with the second or more as

7550

0358	204 Y1(I)=Y2(I)
0229	T1(2)=TPR(4)
0230	CALL DIFEQ (T3,Y1,DY1,M,TPR)
0231	GO TO 8
0232	END .

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